

White Sands Missile Range

Data Review

Otero County, New Mexico

Volatile Organics

SDGs # 9082608, 9090808

Analyses Performed By:

Trace Analysis, Inc.

Lubbock, Texas

Report #10902R

Review Level: Tier II

Project: GP08WSMR00SW OC009

SUMMARY

This data quality assessment summarizes the review of Sample Delivery Group (SDGs) 9082608 and 9090808 for samples collected in association with the White Sands Water Monitoring Site, Otero County, New Mexico. The review was conducted as a Tier II evaluation and included review of data package completeness. Only analytical data associated with constituents of concern were reviewed for this validation. Field documentation was not included in this review. Included with this assessment are the validation annotated sample result sheets, and chain of custody. Analyses were performed on the following samples:

| SDG Number | Sample ID | Lab ID | Matrix | Sample Collection Date | Parent Sample | Analysis | | | | |
|------------|------------------------|--------|--------|------------------------|---------------|----------|------|------------|-----|------|
| | | | | | | VOC | SVOC | HERB /PEST | MET | MISC |
| 9082608 | HLSF-0143-HMW-041-0809 | 207800 | Water | 08/24/2009 | | X | | | | |
| | HLSF-3839-HMW-029-0809 | 207801 | Water | 08/20/2009 | | X | | | | |
| | HLSF-3839-HMW-033-0809 | 207802 | Water | 08/20/2009 | | X | | | | |
| | HLSF-3839-HMW-133-0809 | 207803 | Water | 08/20/2009 | | X | | | | |
| | HLSF-3839-HMW-032-0809 | 207804 | Water | 08/21/2009 | | X | | | | |
| | HLSF-0143-HMW-039-0809 | 207805 | Water | 08/24/2009 | | X | | | | |
| | HLSF-3839-HMW-035-0809 | 207806 | Water | 08/19/2009 | | X | | | | |
| | HLSF-3839-HMW-034-0809 | 207807 | Water | 08/19/2009 | | X | | | | |
| | HLSF-3839-TB-09-000 | 207808 | Water | 08/24/2009 | | X | | | | |
| 9090808 | HLSF-0154-DRW-016-0909 | 209311 | Water | 09/01/2009 | | X | | | | |
| | HLSF-0154-HCF-001-0909 | 209312 | Water | 09/02/2009 | | X | | | | |
| | HLSF-0154-DRW-005-0909 | 209313 | Water | 09/01/2009 | | X | | | | |
| | HLSF-0143-HMW-013-0909 | 209314 | Water | 09/01/2009 | | X | | | | |
| | HLSF-0143-HMW-036-0909 | 209315 | Water | 09/01/2009 | | | | | | |
| | HLSF-0154-DRW-016-0909 | 209316 | Water | 09/01/2009 | MS/MSD | X | | | | |
| | HLSF-0154-DRW-004-0909 | 209317 | Water | 09/02/2009 | | X | | | | |
| | HLSF-0154-TB-09-0002 | 209318 | Water | 09/02/2009 | | X | | | | |

Note: Sample HLSF-0154-DRW-016-0909 was utilized as the Matrix Spike/Matrix Spike Duplicate (MS/MSD).

ANALYTICAL DATA PACKAGE DOCUMENTATION

The table below is the evaluation of the data package completeness.

| Items Reviewed | Reported | | Performance Acceptable | | Not Required |
|---|----------|-----|------------------------|-----|--------------|
| | No | Yes | No | Yes | |
| 1. Sample receipt condition | | X | | X | |
| 2. Requested analyses and sample results | | X | | X | |
| 3. Master tracking list | | X | | X | |
| 4. Methods of analysis | | X | | X | |
| 5. Reporting limits | | X | | X | |
| 6. Sample collection date | | X | | X | |
| 7. Laboratory sample received date | | X | | X | |
| 8. Sample preservation verification (as applicable) | | X | | X | |
| 9. Sample preparation/extraction/analysis dates | | X | | X | |
| 10. Fully executed Chain-of-Custody (COC) form | | X | | X | |
| 11. Narrative summary of QA or sample problems provided | | X | | X | |
| 12. Data Package Completeness and Compliance | | X | | X | |

QA - Quality Assurance

ORGANIC ANALYSIS INTRODUCTION

Analyses were performed according to United States Environmental Protection Agency (USEPA) SW-846 Method 8260B. Data were reviewed in accordance with USEPA National Functional Guidelines of October 1999/January 2005.

The data review process is an evaluation of data on a technical basis rather than a determination of contract compliance. As such, the standards against which the data are being weighed may differ from those specified in the analytical method. It is assumed that the data package represents the best efforts of the laboratory and had already been subjected to adequate and sufficient quality review prior to submission.

During the review process, laboratory qualified and unqualified data are verified against the supporting documentation. Based on this evaluation, qualifier codes may be added, deleted, or modified by the data reviewer. Results are qualified with the following codes in accordance with USEPA National Functional Guidelines:

- Concentration (C) Qualifiers
 - U The compound was analyzed for but not detected. The associated value is the compound quantitation limit.
 - B The compound has been found in the sample as well as its associated blank, its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers
 - E The compound was quantitated above the calibration range.
 - D Concentration is based on a diluted sample analysis.
- Validation Qualifiers
 - J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
 - UJ The compound was not detected above the reported sample quantitation limit. However, the reported limit is approximate and may or may not represent the actual limit of quantitation.
 - JN The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification. The associated numerical value is an estimated concentration only.
 - UB Compound considered non-detect at the listed value due to associated blank contamination.
 - N The analysis indicates the presence of a compound for which there is presumptive evidence to make a tentative identification.
 - R The sample results are rejected.

Two facts should be noted by all data users. First, the "R" flag means that the associated value is unusable. In other words, due to significant quality control (QC) problems, the analysis is invalid and provides no information as to whether the compound is present or not. "R" values should not appear on data tables because they cannot be relied upon, even as a last resort. The second fact to keep in mind is that no compound concentration, even if it has passed all QC tests, is guaranteed to be accurate. Strict QC serves to increase confidence in data but any value potentially contains error.

VOLATILE ORGANIC COMPOUND (VOC) ANALYSES

1. Holding Times

The specified holding times for the following methods are presented in the following table.

| Method | Matrix | Holding Time | Preservation |
|-------------|--------|--|--|
| SW-846 8260 | Water | 14 days from collection to analysis | Cooled @ 4 °C; preserved to a pH of less than 2 s.u. |
| | Soil | 48 hours from collection to extraction and 14 days from extraction to analysis | Cooled @ 4 °C. |

s.u. Standard units

All samples were analyzed within the specified holding time criteria.

2. Blank Contamination

Quality assurance (QA) blanks (i.e., method and rinse blanks) are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Rinse blanks measure contamination of samples during field operations.

A blank action level (BAL) of five times the concentration of a detected compound in an associated blank (common laboratory contaminant compounds are calculated at ten times) is calculated for QA blanks containing concentrations greater than the method detection limit (MDL). The BAL is compared to the associated sample results to determine the appropriate qualification of the sample results, if needed.

All compounds associated with the QA blanks exhibited a concentration less than the MDL, with the exception of the compounds listed in the following table. Sample results associated with QA blank contamination that were greater than the BAL resulted in the removal of the laboratory qualifier (B) of data. Sample results less than the BAL associated with the following sample locations were qualified as listed in the following table.

| Sample Locations | Analytes | Sample Result | Qualification |
|------------------------|------------------------------------|---|--|
| HLSF-0143-HMW-041-0809 | 2-Butanone | Detected sample results >RL and <BAL | "UB" at detected sample concentration |
| HLSF-0154-DRW-005-0909 | Acetone 1,2,4- Trimethylbenzene | | |
| HLSF-0143-HMW-036-0909 | Acetone | | |
| HLSF-0154-DRW-016-0909 | n-Butylbenzene | | |
| HLSF-0154-DRW-004-0909 | Acetone 2-Butanone | | |

RL Reporting limit

3. Surrogates/System Monitoring Compounds

All samples to be analyzed for organic compounds are spiked with surrogate compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. VOC analysis requires that all surrogates associated with the analysis exhibit recoveries within the laboratory-established acceptance limits.

All surrogate recoveries were within control limits.

4. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Analysis

MS/MSD data are used to assess the precision and accuracy of the analytical method. The compounds used to perform the MS/MSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits. The relative percent difference (RPD) between the MS/MSD recoveries must exhibit an RPD within the laboratory-established acceptance limits.

Note: The MS/MSD recovery control limits do not apply for MS/MSD performed on sample locations where the compound concentration detected in the parent sample exceeds the MS/MSD concentration by a factor of four or greater.

Sample locations associated with the MS/MSD exhibiting recoveries outside of the control limits are presented in the following table.

| Sample Locations | Compound | MS Recovery | MSD Recovery |
|------------------------|-----------------------------|--------------|--------------|
| HLSF-0154-DRW-016-0909 | Bromomethane | >UL | >UL |
| | Acetone | <LL but >10% | <LL but >10% |
| | Trans 1,4-Dichloro-2-butene | <LL but >10% | <LL but >10% |
| | Trichloroethene | <LL but >10% | AC |
| | 2-Chloroethyl vinyl ether | < 10% | < 10% |
| | Tetrachloroethene | <LL but >10% | <LL but >10% |

AC Acceptable

The criteria used to evaluate the MS/MSD recoveries are presented in the following table. In the case of an MS/MSD deviation, the sample results are qualified as documented in the table below.

| Control Limit | Sample Result | Qualification |
|---|---------------|---------------|
| > the upper control limit (UL) | Non-detect | No Action |
| | Detect | J |
| < the lower control limit (LL) but > 10% | Non-detect | UJ |
| | Detect | J |
| < 10% | Non-detect | R |
| | Detect | J |
| Parent sample concentration > four times the MS/MSD spiking solution concentration (D). | Detect | No Action |
| | Non-detect | |

5. Laboratory Control Sample/Laboratory Control Sample Duplicate (LCS/LCSD) Analysis

The LCS/LCSD analysis is used to assess the precision and accuracy of the analytical method independent of matrix interferences. The compounds associated with the LCS/LCSD analysis must exhibit a percent recovery within the laboratory-established acceptance limits.

Sample locations associated with LCS/LCSD analysis exhibiting recoveries outside of the control limits presented in the following table.

| Sample Locations | Compound | LCS Recovery | LCSD Recovery |
|---|---------------------------|--|----------------|
| SDG: 9082608 HLSF-0143-HMW-041-0809 HLSF-3839-HMW-029-0809 HLSF-3839-HMW-033-0809 HLSF-3839-HMW-133-0809 HLSF-3839-HMW-032-0809 HLSF-0143-HMW-039-0809 HLSF-3839-HMW-035-0809 HLSF-3839-HMW-034-0809 HLSF-3839-TB-09-000 | 1,2-Dichloroethane | < LL but > 10% < LL but > 10% < LL but > 10% < LL but > 10% < LL but > 10% < LL but > 10% < LL but > 10% | < LL but > 10% |
| | 1,1,1-Trichloroethane | | |
| | Carbon Tetrachloride | | |
| | 1,1,1,2-Tetrachloroethane | | |
| | Bromoform | | |
| | 1,4-Dichlorobenzene | | |
| | Dibromomethane | AC | < LL but > 10% |
| SDG: 9090808 HLSF-0154-DRW-016-0909 HLSF-0154-HCF-001-0909 HLSF-0154-DRW-005-0909 HLSF-0143-HMW-013-0909 HLSF-0143-HMW-036-0909 HLSF-0154-DRW-016-0909 HLSF-0154-DRW-004-0909 HLSF-0154-TB-09-0002 | Acetone | < LL but > 10% | AC |
| | Terachloroethene | < LL but > 10% | AC |

The criteria used to evaluate the LCS/LCSD recoveries are presented in the following table. In the case of an LCS/LCSD deviation, the sample results are qualified as documented in the table below.

| Control Limit | Sample Result | Qualification |
|--|---------------|---------------|
| > the upper control limit (UL) | Non-detect | No Action |
| | Detect | J |
| < the lower control limit (LL) but > 10% | Non-detect | UJ |
| | Detect | J |
| < 10% | Non-detect | R |
| | Detect | J |

Note: Sample results were not qualified as rejected (R) due to the deviations listed above.

6. Field Duplicate Analysis

Field duplicate analysis is used to assess the precision and accuracy of the field sampling procedures and analytical method. A control limit of 50% for water matrices is applied to the RPD between the parent sample and the field duplicate. In the instance when the parent and/or duplicate sample concentrations are less than or equal to 5 times the RL, a control limit of two times the RL is applied for water matrices.

A field duplicate was not collected with a sample location associated with this SDG.

7. Compound Identification

Compounds are identified on the GC/MS by laboratory personnel using the analytes relative retention time and ion spectra.

8. System Performance and Overall Assessment

Overall system performance was noted by the laboratory in the case narrative as acceptable. Other than for those deviations specifically mentioned in this review, the overall data quality is within the guidelines specified in the method.

DATA VALIDATION CHECKLIST FOR VOCs

| VOCs: SW-846 8260 | Reported | | Performance Acceptable | | Not Required |
|--|----------|-----|------------------------|-----|--------------|
| | No | Yes | No | Yes | |
| GAS CHROMATOGRAPHY/MASS SPECTROMETRY (GC/MS) | | | | | |
| Tier II Validation | | | | | |
| Holding times | | X | | X | |
| Reporting limits (units) | | X | | X | |
| Blanks | | | | | |
| A. Method blanks | | X | X | | |
| B. Equipment blanks | | | | | X |
| C. Trip blanks | | X | X | X | |
| Laboratory Control Sample (LCS) | | X | X | | |
| Laboratory Control Sample Duplicate(LCSD) | | X | X | | |
| LCS/LCSD Precision (RPD) | | X | | X | |
| Matrix Spike (MS) | | X | X | | |
| Matrix Spike Duplicate(MSD) | | X | X | | |
| MS/MSD Precision (RPD) | | X | | X | |
| Field/Lab Duplicate (RPD) | | | | | X |
| Surrogate Spike Recoveries | | X | | X | |
| Dilution Factor | | X | | X | |
| Moisture Content | | | | | X |

%RSD Relative standard deviation

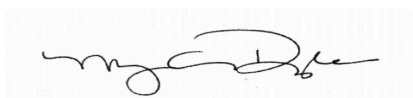
%R Percent recovery

RPD Relative percent difference

%D Percent difference

VALIDATION PERFORMED BY: Mary Ann Doyle

SIGNATURE:

A handwritten signature in black ink, appearing to read 'Mary Ann Doyle', is written over a light gray rectangular background. The signature is fluid and cursive.

DATE: October 15, 2009

PEER REVIEW: Dennis Capria

DATE: October 19, 2009

**CHAIN OF CUSTODY/
CORRECTED SAMPLE ANALYSIS DATA SHEETS**

Analytical Report

Sample: 207800 - HLSF-0143-HMW-041-0809

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 62999

Prep Batch: 53764

Analytical Method: S 8260B

Date Analyzed: 2009-08-27

Sample Preparation: 2009-08-27

Prep Method: S 5030B

Analyzed By: KB

Prepared By: KB

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|---------------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | ^U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | ^{1 U} | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | ^U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | ^U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | ^U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |
| Chloroethane | ^U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | ^{2 U} | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | ^U | <1.75 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | ^U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | ^U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | ^U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | ^{UB 3 3 3} | 0.880 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | ^U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | ^U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans 1,4-Dichloro-2-butene | ^{4 U} | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | | 3.79 | 3.79 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | ^U | <0.450 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | ^U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | ^U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | ^J | 0.890 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| cis-1,2-Dichloroethene | ^U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | ^U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | ^{5 U J} | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | | 1.14 | 1.14 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | ^{6 U J} | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | ^U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | ^U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| Carbon Tetrachloride | ^{7 U J} | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | ^U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | | 118 | 118 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | ^{8 U J} | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |

continued ...

¹ Concentration biased low.² Concentration biased low.³ Concentration biased high.⁴ Concentration biased low.⁵ Concentration biased low.⁶ Concentration biased low.⁷ Concentration biased low.⁸ Concentration biased low.

sample 207800 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|-----------------------------|---------------------------------|-----------------|-----------------|-----------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromodichloromethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| cis-1,3-Dichloropropene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichloropropane | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Tetrachloroethene (PCE) | ⁹ U | 0.400 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,1,2-Tetrachloroethane | ¹⁰ U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | ¹¹ U ¹² J | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| 1,2,3-Trichloropropane | U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | ¹² U ¹³ J | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | ¹³ U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

⁹Concentration biased low.¹⁰Concentration biased low.¹¹Concentration biased low.¹²Concentration biased low.¹³Concentration biased low.

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 49.8 | µg/L | 1 | 50.0 | 100 | 75.3 - 131 |
| Toluene-d8 | | 54.1 | µg/L | 1 | 50.0 | 108 | 91.4 - 112 |
| 4-Bromofluorobenzene (4-BFB) | | 50.1 | µg/L | 1 | 50.0 | 100 | 83.8 - 108 |

Sample: 207801 - HLSF-3839-HMW-029-0809

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 62999

Prep Batch: 53764

Analytical Method: S 8260B

Date Analyzed: 2009-08-27

Sample Preparation: 2009-08-27

Prep Method: S 5030B

Analyzed By: KB

Prepared By: KB

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|---------------------------------|-------|--------------|--------------|--------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | N U | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |
| Chloroethane | U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | N U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | U | <1.75 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | N U | <0.810 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans 1,4-Dichloro-2-butene | N U | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | U | <0.450 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| cis-1,2-Dichloroethene | U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | N U J | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | N U J | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |

*continued ...*¹⁴Concentration biased low.¹⁵Concentration biased low.¹⁶Concentration biased high.¹⁷Concentration biased low.¹⁸Concentration biased low.¹⁹Concentration biased low.

sample 207801 continued ...

| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|------------------------------------|-----------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| Carbon Tetrachloride | 20 U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | 21 U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Bromodichloromethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| cis-1,3-Dichloropropene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichloropropane | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Tetrachloroethene (PCE) | 22 U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | 23 U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,1,2-Tetrachloroethane | 24 U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | 25 U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| 1,2,3-Trichloropropane | U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | 26 U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | 27 U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | 28 U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |

continued ...

²⁰Concentration biased low.²¹Concentration biased low.²²Concentration biased low.²³Concentration biased low.²⁴Concentration biased low.²⁵Concentration biased low.²⁶Concentration biased low.

sample 207801 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------|------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 50.0 | µg/L | 1 | 50.0 | 100 | 75.3 - 131 |
| Toluene-d8 | | 52.5 | µg/L | 1 | 50.0 | 105 | 91.4 - 112 |
| 4-Bromofluorobenzene (4-BFB) | | 49.2 | µg/L | 1 | 50.0 | 98 | 83.8 - 108 |

Sample: 207802 - HLSF-3839-HMW-033-0809

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 62999

Prep Batch: 53764

Analytical Method: S 8260B

Date Analyzed: 2009-08-27

Sample Preparation: 2009-08-27

Prep Method: S 5030B

Analyzed By: KB

Prepared By: KB

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|---------------------------------|-----------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | 27 U | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |
| Chloroethane | U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | 28 U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | U | <1.75 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | 29 U | <0.810 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans 1,4-Dichloro-2-butene | 30 U | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | | 4.15 | 4.15 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | U | <0.450 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |

continued ...

²⁷Concentration biased low.²⁸Concentration biased low.²⁹Concentration biased high.³⁰Concentration biased low.

sample 207802 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|------------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| cis-1,2-Dichloroethene | ³¹ U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | ³² U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | ³³ UJ | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | ³⁴ UJ | 1.36 | 1.36 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | ³⁵ UJ | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | ³⁶ U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | ³⁷ U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| Carbon Tetrachloride | ³⁸ UJ | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | ³⁹ U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | ⁴⁰ UJ | 4.75 | 4.75 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | ⁴¹ UJ | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Bromodichloromethane | ⁴² U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | ⁴³ U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| cis-1,3-Dichloropropene | ⁴⁴ U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | ⁴⁵ U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | ⁴⁶ U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | ⁴⁷ U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichloropropane | ⁴⁸ U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | ⁴⁹ U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | ⁵⁰ U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Tetrachloroethene (PCE) | ⁵¹ J | 0.430 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | ⁵² U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,1,2-Tetrachloroethane | ⁵³ U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | ⁵⁴ U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | ⁵⁵ U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | ⁵⁶ UJ | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | ⁵⁷ U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | ⁵⁸ U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | ⁵⁹ U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | ⁶⁰ U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| 1,2,3-Trichloropropane | ⁶¹ U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | ⁶² U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | ⁶³ U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | ⁶⁴ U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | ⁶⁵ U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | ⁶⁶ U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | ⁶⁷ U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | ⁶⁸ UJ | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |

continued ...

³¹Concentration biased low.³²Concentration biased low.³³Concentration biased low.³⁴Concentration biased low.³⁵Concentration biased low.³⁶Concentration biased low.³⁷Concentration biased low.³⁸Concentration biased low.

sample 207802 continued ...

| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|-----------------------------|-----------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | 39 U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|-----------------|---------------------|--------------------|
| Dibromofluoromethane | | 49.2 | µg/L | 1 | 50.0 | 98 | 75.3 - 131 |
| Toluene-d8 | | 52.8 | µg/L | 1 | 50.0 | 106 | 91.4 - 112 |
| 4-Bromofluorobenzene (4-BFB) | | 48.9 | µg/L | 1 | 50.0 | 98 | 83.8 - 108 |

Sample: 207803 - HLSF-3839-HMW-133-0809

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 62999

Prep Batch: 53764

Analytical Method: S 8260B

Date Analyzed: 2009-08-27

Sample Preparation: 2009-08-27

Prep Method: S 5030B

Analyzed By: KB

Prepared By: KB

| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|---------------------------------|-----------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| Bromochloromethane | U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | 40 U | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |
| Chloroethane | U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | 41 U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | U | <1.75 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | 42 U | <0.810 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |

continued ...

³⁹Concentration biased low.⁴⁰Concentration biased low.⁴¹Concentration biased low.⁴²Concentration biased high.

sample 207803 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|-------------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| 4-Methyl-2-pentanone (MIBK) | U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans 1,4-Dichloro-2-butene | 43 U | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | J | 4.26 | 4.26 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | U | <0.450 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| cis-1,2-Dichloroethene | U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | 44 U J | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | | 1.35 | 1.35 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | 45 U J | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| Carbon Tetrachloride | 46 U J | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | | 4.85 | 4.85 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | 47 U J | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Bromodichloromethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| cis-1,3-Dichloropropene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichloropropane | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Tetrachloroethene (PCE) | 48 U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,1,2-Tetrachloroethane | 49 U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | 50 U J | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |

continued ...

⁴³Concentration biased low.⁴⁴Concentration biased low.⁴⁵Concentration biased low.⁴⁶Concentration biased low.⁴⁷Concentration biased low.⁴⁸Concentration biased low.⁴⁹Concentration biased low.⁵⁰Concentration biased low.

sample 207803 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|-----------------------------|------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| 1,2,3-Trichloropropane | U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 51.5 | µg/L | 1 | 50.0 | 103 | 75.3 - 131 |
| Toluene-d8 | | 52.9 | µg/L | 1 | 50.0 | 106 | 91.4 - 112 |
| 4-Bromofluorobenzene (4-BFB) | | 49.2 | µg/L | 1 | 50.0 | 98 | 83.8 - 108 |

Sample: 207804 - HLSF-3839-HMW-032-0809

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 62999

Prep Batch: 53764

Analytical Method: S 8260B

Date Analyzed: 2009-08-27

Sample Preparation: 2009-08-27

Prep Method: S 5030B

Analyzed By: KB

Prepared By: KB

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|---------------------------------|------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | U | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |

continued ...

⁵¹ Concentration biased low.⁵² Concentration biased low.⁵³ Concentration biased low.

sample 207804 continued ...

| Parameter | Flag | SDL Based Result | SQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|------------------------------------|-----------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| Chloroethane | U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | ⁵⁴ U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | U | <1.75 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | 55 U | <0.810 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans 1,4-Dichloro-2-butene | 56 U | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | U | <0.450 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| cis-1,2-Dichloroethene | U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | 57 U | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | 58 U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| Carbon Tetrachloride | 59 U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | 60 U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Bromodichloromethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| cis-1,3-Dichloropropene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichloropropane | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Tetrachloroethene (PCE) | 61 U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |

continued ...

⁵⁴Concentration biased low.⁵⁵Concentration biased high.⁵⁶Concentration biased low.⁵⁷Concentration biased low.⁵⁸Concentration biased low.⁵⁹Concentration biased low.⁶⁰Concentration biased low.⁶¹Concentration biased low.

sample 207804 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|-----------------------------|-----------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| 1,1,1,2-Tetrachloroethane | 62 U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | 63 U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| 1,2,3-Trichloropropane | U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | 64 U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | 65 U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 51.4 | µg/L | 1 | 50.0 | 103 | 75.3 - 131 |
| Toluene-d8 | | 52.6 | µg/L | 1 | 50.0 | 105 | 91.4 - 112 |
| 4-Bromofluorobenzene (4-BFB) | | 49.8 | µg/L | 1 | 50.0 | 100 | 83.8 - 108 |

Sample: 207805 - HLSF-0143-HMW-039-0809

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 62999

Prep Batch: 53764

Analytical Method: S 8260B

Date Analyzed: 2009-08-27

Sample Preparation: 2009-08-27

Prep Method: S 5030B

Analyzed By: KB

Prepared By: KB

⁶²Concentration biased low.⁶³Concentration biased low.⁶⁴Concentration biased low.⁶⁵Concentration biased low.

| Parameter | Flag | SDL Based Result | SQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|------------------------------------|--------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| Bromochloromethane | U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | U | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |
| Chloroethane | U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | U | <1.75 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | U | <0.810 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans-1,4-Dichloro-2-butene | U | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | U | 1.53 | 1.53 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | U | <0.450 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| cis-1,2-Dichloroethene | U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | U | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | U | 0.440 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| Carbon Tetrachloride | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | U | 44.0 | 44.0 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Bromodichloromethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| cis-1,3-Dichloropropene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichloropropane | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |

continued ...

⁶⁶Concentration biased low.⁶⁷Concentration biased low.⁶⁸Concentration biased high.⁶⁹Concentration biased low.⁷⁰Concentration biased low.⁷¹Concentration biased low.⁷²Concentration biased low.⁷³Concentration biased low.

sample 207805 continued ...

| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|-----------------------------|------------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| Tetrachloroethene (PCE) | N U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,1,2-Tetrachloroethane | N U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | N U J | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| 1,2,3-Trichloropropane | U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | N U J | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | N U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | N U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|-----------------|---------------------|--------------------|
| Dibromofluoromethane | | 50.3 | µg/L | 1 | 50.0 | 101 | 75.3 - 131 |
| Toluene-d8 | | 52.4 | µg/L | 1 | 50.0 | 105 | 91.4 - 112 |
| 4-Bromofluorobenzene (4-BFB) | | 49.7 | µg/L | 1 | 50.0 | 99 | 83.8 - 108 |

Sample: 207806 - HLSF-3839-HMW-035-0809

Laboratory: Lubbock

Analysis: Volatiles

Analytical Method: S 8260B

Prep Method: S 5030B

⁷⁴Concentration biased low.⁷⁵Concentration biased low.⁷⁶Concentration biased low.⁷⁷Concentration biased low.⁷⁸Concentration biased low.

QC Batch: 62999
Prep Batch: 53764Date Analyzed: 2009-08-27
Sample Preparation: 2009-08-27Analyzed By: KB
Prepared By: KB

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|------------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | ⁷⁹ U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | ⁷⁹ U | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |
| Chloroethane | ⁸⁰ U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | ⁸⁰ U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | U | <1.75 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | ⁸¹ U | <0.810 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans-1,4-Dichloro-2-butene | ⁸² U | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | U | <0.450 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| cis-1,2-Dichloroethene | U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | ⁸³ UJ | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | ⁸⁴ UJ | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| Carbon Tetrachloride | ⁸⁵ UJ | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | ⁸⁶ UJ | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Bromodichloromethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| cis-1,3-Dichloropropene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |

continued ...

⁷⁹Concentration biased low.⁸⁰Concentration biased low.⁸¹Concentration biased high.⁸²Concentration biased low.⁸³Concentration biased low.⁸⁴Concentration biased low.⁸⁵Concentration biased low.⁸⁶Concentration biased low.

sample 207806 continued ...

| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|-----------------------------|-----------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| 1,3-Dichloropropane | ⁸⁷ U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Tetrachloroethene (PCE) | ⁸⁸ U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | ⁸⁸ U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,1,2-Tetrachloroethane | ⁸⁸ U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | ⁸⁹ U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| 1,2,3-Trichloropropane | U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | ⁹⁰ U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | ⁹¹ U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|-----------------|---------------------|--------------------|
| Dibromofluoromethane | | 51.5 | µg/L | 1 | 50.0 | 103 | 75.3 - 131 |
| Toluene-d8 | | 52.4 | µg/L | 1 | 50.0 | 105 | 91.4 - 112 |
| 4-Bromofluorobenzene (4-BFB) | | 50.4 | µg/L | 1 | 50.0 | 101 | 83.8 - 108 |

⁸⁷Concentration biased low.⁸⁸Concentration biased low.⁸⁹Concentration biased low.⁹⁰Concentration biased low.⁹¹Concentration biased low.

Sample: 207807 - HLSF-3839-HMW-034-0809

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 62999

Prep Batch: 53764

Analytical Method: S 8260B

Date Analyzed: 2009-08-27

Sample Preparation: 2009-08-27

Prep Method: S 5030B

Analyzed By: KB

Prepared By: KB

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|-------------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | U | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |
| Chloroethane | U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | U | <1.75 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | 92 U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | 98 U | <0.810 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans 1,4-Dichloro-2-butene | U | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | U | <0.450 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| cis-1,2-Dichloroethene | U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | 94 U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | 95 U J | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | 96 U J | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| Carbon Tetrachloride | 97 U J | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | 98 U J | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Bromodichloromethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | 99 U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |

continued ...

⁹²Concentration biased high.⁹³Concentration biased high.⁹⁴Concentration biased low.⁹⁵Concentration biased low.⁹⁶Concentration biased low.⁹⁷Concentration biased low.⁹⁸Concentration biased low.⁹⁹Concentration biased high.

sample 207807 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|-----------------------------|------------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| cis-1,3-Dichloropropene | ¹⁰⁰ U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichloropropane | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Tetrachloroethene (PCE) | ¹⁰¹ U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,1,2-Tetrachloroethane | ¹⁰² U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | ¹⁰³ U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| 1,2,3-Trichloropropane | U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | ¹⁰⁴ U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|----------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 53.9 | µg/L | 1 | 50.0 | 108 | 75.3 - 131 |
| Toluene-d8 | | 52.7 | µg/L | 1 | 50.0 | 105 | 91.4 - 112 |

continued ...

¹⁰⁰Concentration biased high.¹⁰¹Concentration biased high.¹⁰²Concentration biased low.¹⁰³Concentration biased low.¹⁰⁴Concentration biased low.

sample continued ...

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| 4-Bromofluorobenzene (4-BFB) | | 51.4 | µg/L | 1 | 50.0 | 103 | 83.8 - 108 |

Sample: 207808 - HLSF-3839-TB-09-000

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 62999

Prep Batch: 53764

Analytical Method: S 8260B

Date Analyzed: 2009-08-27

Sample Preparation: 2009-08-27

Prep Method: S 5030B

Analyzed By: KB

Prepared By: KB

| Parameter | Flag | SDL MQL Method | | | Units | Dilution | MQL | | MDL |
|---------------------------------|------------------|----------------|-------|--------|-------|----------|-------|--------------|--------------|
| | | Based | Based | Blank | | | SDL | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | U | <0.370 | <1.00 | <0.370 | µg/L | 1 | 0.370 | 1 | 0.37 |
| Dichlorodifluoromethane | 105 U | <0.450 | <1.00 | <0.450 | µg/L | 1 | 0.450 | 1 | 0.45 |
| Chloromethane (methyl chloride) | U | <0.590 | <1.00 | <0.590 | µg/L | 1 | 0.590 | 1 | 0.59 |
| Vinyl Chloride | U | <0.690 | <1.00 | <0.690 | µg/L | 1 | 0.690 | 1 | 0.69 |
| Bromomethane (methyl bromide) | U | <0.750 | <5.00 | <0.750 | µg/L | 1 | 0.750 | 5 | 0.75 |
| Chloroethane | U | <0.570 | <1.00 | <0.570 | µg/L | 1 | 0.570 | 1 | 0.57 |
| Trichlorofluoromethane | 106 U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Acetone | JB | 3.08 | <10.0 | 8.53 | µg/L | 1 | 1.75 | 10 | 1.75 |
| Iodomethane (methyl iodide) | U | <0.320 | <5.00 | <0.320 | µg/L | 1 | 0.320 | 5 | 0.32 |
| Carbon Disulfide | U | <0.250 | <1.00 | <0.250 | µg/L | 1 | 0.250 | 1 | 0.25 |
| Acrylonitrile | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 2-Butanone (MEK) | 107 U | <0.810 | <5.00 | 11.3 | µg/L | 1 | 0.810 | 5 | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.790 | <5.00 | <0.790 | µg/L | 1 | 0.790 | 5 | 0.79 |
| 2-Hexanone | U | <0.510 | <5.00 | <0.510 | µg/L | 1 | 0.510 | 5 | 0.51 |
| trans 1,4-Dichloro-2-butene | 108 U | <0.490 | <10.0 | <0.490 | µg/L | 1 | 0.490 | 10 | 0.49 |
| 1,1-Dichloroethene | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| Methylene chloride | J | 1.92 | <5.00 | <0.450 | µg/L | 1 | 0.450 | 5 | 0.45 |
| MTBE | U | <0.400 | <1.00 | <0.400 | µg/L | 1 | 0.400 | 1 | 0.4 |
| trans-1,2-Dichloroethene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 1,1-Dichloroethane | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| cis-1,2-Dichloroethene | U | <0.200 | <1.00 | <0.200 | µg/L | 1 | 0.200 | 1 | 0.2 |
| 2,2-Dichloropropane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 1,2-Dichloroethane (EDC) | 109 U | <0.350 | <1.00 | <0.350 | µg/L | 1 | 0.350 | 1 | 0.35 |
| Chloroform | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,1-Trichloroethane | 110 U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| 1,1-Dichloropropene | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Benzene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |

continued ...

¹⁰⁵Concentration biased low.¹⁰⁶Concentration biased low.¹⁰⁷Concentration biased high.¹⁰⁸Concentration biased low.¹⁰⁹Concentration biased low.¹¹⁰Concentration biased low.

sample 207808 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|------------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Carbon Tetrachloride | 111 U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dichloropropane | U | <0.360 | <1.00 | <0.360 | µg/L | 1 | 0.360 | 1 | 0.36 |
| Trichloroethene (TCE) | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| Dibromomethane (methylene bromide) | 112 U | <0.470 | <1.00 | <0.470 | µg/L | 1 | 0.470 | 1 | 0.47 |
| Bromodichloromethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 2-Chloroethyl vinyl ether | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |
| cis-1,3-Dichloropropene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| trans-1,3-Dichloropropene | U | <0.380 | <1.00 | <0.380 | µg/L | 1 | 0.380 | 1 | 0.38 |
| Toluene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| 1,1,2-Trichloroethane | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichloropropane | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| Dibromochloromethane | U | <0.320 | <1.00 | <0.320 | µg/L | 1 | 0.320 | 1 | 0.32 |
| 1,2-Dibromoethane (EDB) | U | <0.340 | <1.00 | <0.340 | µg/L | 1 | 0.340 | 1 | 0.34 |
| Tetrachloroethene (PCE) | 113 U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| Chlorobenzene | 114 U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,1,2-Tetrachloroethane | 115 U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Ethylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| m,p-Xylene | U | <0.540 | <1.00 | <0.540 | µg/L | 1 | 0.540 | 1 | 0.54 |
| Bromoform | 116 U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| Styrene | U | <0.210 | <1.00 | <0.210 | µg/L | 1 | 0.210 | 1 | 0.21 |
| o-Xylene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| 1,1,2,2-Tetrachloroethane | U | <0.420 | <1.00 | <0.420 | µg/L | 1 | 0.420 | 1 | 0.42 |
| 2-Chlorotoluene | U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| 1,2,3-Trichloropropane | U | <0.430 | <1.00 | <0.430 | µg/L | 1 | 0.430 | 1 | 0.43 |
| Isopropylbenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| Bromobenzene | U | <0.260 | <1.00 | <0.260 | µg/L | 1 | 0.260 | 1 | 0.26 |
| n-Propylbenzene | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| 1,3,5-Trimethylbenzene | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| tert-Butylbenzene | U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2,4-Trimethylbenzene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,4-Dichlorobenzene (para) | 117 U | <0.240 | <1.00 | <0.240 | µg/L | 1 | 0.240 | 1 | 0.24 |
| sec-Butylbenzene | U | <0.280 | <1.00 | <0.280 | µg/L | 1 | 0.280 | 1 | 0.28 |
| 1,3-Dichlorobenzene (meta) | U | <0.310 | <1.00 | <0.310 | µg/L | 1 | 0.310 | 1 | 0.31 |
| p-Isopropyltoluene | U | <0.330 | <1.00 | <0.330 | µg/L | 1 | 0.330 | 1 | 0.33 |
| 4-Chlorotoluene | U | <0.290 | <1.00 | <0.290 | µg/L | 1 | 0.290 | 1 | 0.29 |
| 1,2-Dichlorobenzene (ortho) | U | <0.270 | <1.00 | <0.270 | µg/L | 1 | 0.270 | 1 | 0.27 |
| n-Butylbenzene | 118 U | <0.300 | <1.00 | <0.300 | µg/L | 1 | 0.300 | 1 | 0.3 |
| 1,2-Dibromo-3-chloropropane | 119 U | <0.680 | <5.00 | <0.680 | µg/L | 1 | 0.680 | 5 | 0.68 |
| 1,2,3-Trichlorobenzene | U | <0.330 | <5.00 | <0.330 | µg/L | 1 | 0.330 | 5 | 0.33 |

continued ...

¹¹¹Concentration biased low.¹¹²Concentration biased low.¹¹³Concentration biased low.¹¹⁴Concentration biased low.¹¹⁵Concentration biased low.¹¹⁶Concentration biased low.¹¹⁷Concentration biased low.

sample 207808 continued ...

| Parameter | Flag | SDL | SQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------|------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| 1,2,4-Trichlorobenzene | U | <0.340 | <5.00 | <0.340 | µg/L | 1 | 0.340 | 5 | 0.34 |
| Naphthalene | U | <0.280 | <5.00 | <0.280 | µg/L | 1 | 0.280 | 5 | 0.28 |
| Hexachlorobutadiene | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 49.2 | µg/L | 1 | 50.0 | 98 | 75.3 - 131 |
| Toluene-d8 | | 53.8 | µg/L | 1 | 50.0 | 108 | 91.4 - 112 |
| 4-Bromofluorobenzene (4-BFB) | | 50.6 | µg/L | 1 | 50.0 | 101 | 83.8 - 108 |

Method Blank (1)

QC Batch: 62999

Date Analyzed: 2009-08-27

Analyzed By: KB

Prep Batch: 53764

QC Preparation: 2009-08-27

Prepared By: KB

| Parameter | Flag | Result | Units | Reporting Limits |
|---------------------------------|------|--------|-------|------------------|
| Bromochloromethane | | <0.370 | µg/L | 0.37 |
| Dichlorodifluoromethane | | <0.450 | µg/L | 0.45 |
| Chloromethane (methyl chloride) | | <0.590 | µg/L | 0.59 |
| Vinyl Chloride | | <0.690 | µg/L | 0.69 |
| Bromomethane (methyl bromide) | | <0.750 | µg/L | 0.75 |
| Chloroethane | | <0.570 | µg/L | 0.57 |
| Trichlorofluoromethane | | <0.470 | µg/L | 0.47 |
| Acetone | | 8.53 | µg/L | 1.75 |
| Iodomethane (methyl iodide) | | <0.320 | µg/L | 0.32 |
| Carbon Disulfide | | <0.250 | µg/L | 0.25 |
| Acrylonitrile | | <0.320 | µg/L | 0.32 |
| 2-Butanone (MEK) | | 11.3 | µg/L | 0.81 |
| 4-Methyl-2-pentanone (MIBK) | | <0.790 | µg/L | 0.79 |
| 2-Hexanone | | <0.510 | µg/L | 0.51 |
| trans 1,4-Dichloro-2-butene | | <0.490 | µg/L | 0.49 |
| 1,1-Dichloroethene | | <0.400 | µg/L | 0.4 |
| Methylene chloride | | <0.450 | µg/L | 0.45 |
| MTBE | | <0.400 | µg/L | 0.4 |
| trans-1,2-Dichloroethene | | <0.330 | µg/L | 0.33 |
| 1,1-Dichloroethane | | <0.290 | µg/L | 0.29 |
| cis-1,2-Dichloroethene | | <0.200 | µg/L | 0.2 |
| 2,2-Dichloropropane | | <0.420 | µg/L | 0.42 |
| 1,2-Dichloroethane (EDC) | | <0.350 | µg/L | 0.35 |
| Chloroform | | <0.270 | µg/L | 0.27 |
| 1,1,1-Trichloroethane | | <0.230 | µg/L | 0.23 |
| 1,1-Dichloropropene | | <0.340 | µg/L | 0.34 |
| Benzene | | <0.240 | µg/L | 0.24 |

continued ...

1000 S. JEFFERSON BLVD STE 100
 LEBANON, NH 03601
 603-407-1500
 603-407-1501 F

CHAIN OF CUSTODY RECORD

| PROJECT NO. | | PROJECT NAME | | ANALYSIS REQUIRED | |
|-------------|------|-------------------------------|---------|-------------------|-------------------|
| 9032408 | | HESTF Semi-Annual Groundwater | | VOCs | |
| DATE | TIME | SAMPLE ID | ANALYST | LAB NO. | NO. OF CONTAINERS |
| 8-24-09 | 1015 | HLSF-0143-HMW-041-0809 | Water | 307800 | 3 |
| 8-20-09 | 1354 | HLSF-3839-HMW-039-0809 | Water | 801 | 3 |
| 8-20-09 | 0958 | HLSF-3839-HMW-033-0809 | Water | 802 | 3 |
| 8-20-09 | 0958 | HLSF-3839-HMW-133-0809 | Water | 803 | 3 |
| 8-21-09 | 1340 | HLSF-3839-HMW-032-0809 | Water | 804 | 3 |
| 8-24-09 | 1310 | HLSF-0143-HMW-039-0809 | Water | 805 | 3 |
| 8-19-09 | 1305 | HLSF-3839-HMW-035-0809 | Water | 806 | 3 |
| 8-19-09 | 1017 | HLSF-3839-HMW-034-0809 | Water | 807 | 3 |
| 8-24-09 | 1310 | HLSF-3839-TB-09-000 | Water | 808 | 3 |

| 1. PREPARED BY (SIGNATURE) | 2. PREPARED BY (SIGNATURE) | 3. PREPARED BY (SIGNATURE) |
|-------------------------------------|----------------------------|---------------------------------|
| Bradley T. Davis 8-25-09 1300 | | Carol Fox 8-26-09 8:30 AM |

| DATE | TIME | ANALYST | LAB NO. | NO. OF CONTAINERS |
|---------|------|---------|---------|-------------------|
| 8-24-09 | 1015 | Water | 307800 | 3 |
| 8-20-09 | 1354 | Water | 801 | 3 |
| 8-20-09 | 0958 | Water | 802 | 3 |
| 8-20-09 | 0958 | Water | 803 | 3 |
| 8-21-09 | 1340 | Water | 804 | 3 |
| 8-24-09 | 1310 | Water | 805 | 3 |
| 8-19-09 | 1305 | Water | 806 | 3 |
| 8-19-09 | 1017 | Water | 807 | 3 |
| 8-24-09 | 1310 | Water | 808 | 3 |

10902R (2)

Analytical Report

Sample: 209311 - HLSF-0154-DRW-016-0909

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 63556

Prep Batch: 54254

Analytical Method: S 8260B

Date Analyzed: 2009-09-14

Sample Preparation: 2009-09-14

Prep Method: S 5030B

Analyzed By: JG

Prepared By: JG

| Parameter | Flag | SDL Based Result | SQL Based Result | Method Blank Result | Units | Dilution | SDL | SQL (Unadjusted) | MDL (Unadjusted) |
|------------------------------------|------|------------------------|------------------------|---------------------------|-------|----------|--------|---------------------|---------------------|
| Bromochloromethane | U | <0.169 | <1.00 | <0.169 | µg/L | 1 | 0.169 | 1 | 0.169 |
| Dichlorodifluoromethane | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Chloromethane (methyl chloride) | U | <0.164 | <1.00 | <0.164 | µg/L | 1 | 0.164 | 1 | 0.164 |
| Vinyl Chloride | U | <0.110 | <1.00 | <0.110 | µg/L | 1 | 0.110 | 1 | 0.11 |
| Bromomethane (methyl bromide) | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |
| Chloroethane | U | <0.232 | <1.00 | <0.232 | µg/L | 1 | 0.232 | 1 | 0.232 |
| Trichlorofluoromethane | J | 0.394 | <1.00 | <0.149 | µg/L | 1 | 0.149 | 1 | 0.149 |
| Acetone | U J | <0.893 | <10.0 | 2.05 | µg/L | 1 | 0.893 | 10 | 0.893 |
| Iodomethane (methyl iodide) | U | <0.104 | <5.00 | <0.104 | µg/L | 1 | 0.104 | 5 | 0.104 |
| Carbon Disulfide | U | <0.120 | <1.00 | <0.120 | µg/L | 1 | 0.120 | 1 | 0.12 |
| Acrylonitrile | U | <0.189 | <1.00 | <0.189 | µg/L | 1 | 0.189 | 1 | 0.189 |
| 2-Butanone (MEK) | U | <0.218 | <5.00 | 0.596 | µg/L | 1 | 0.218 | 5 | 0.218 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.554 | <5.00 | <0.554 | µg/L | 1 | 0.554 | 5 | 0.554 |
| 2-Hexanone | U | <0.863 | <5.00 | <0.863 | µg/L | 1 | 0.863 | 5 | 0.863 |
| trans 1,4-Dichloro-2-butene | U J | <0.607 | <10.0 | <0.607 | µg/L | 1 | 0.607 | 10 | 0.607 |
| 1,1-Dichloroethene | U | 5.84 | 5.84 | <0.102 | µg/L | 1 | 0.102 | 1 | 0.102 |
| Methylene chloride | U | <0.270 | <5.00 | <0.270 | µg/L | 1 | 0.270 | 5 | 0.27 |
| MTBE | J | 0.289 | <1.00 | <0.122 | µg/L | 1 | 0.122 | 1 | 0.122 |
| trans-1,2-Dichloroethene | U | <0.168 | <1.00 | <0.168 | µg/L | 1 | 0.168 | 1 | 0.168 |
| 1,1-Dichloroethane | U | 1.32 | 1.32 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| cis-1,2-Dichloroethene | U | <0.151 | <1.00 | <0.151 | µg/L | 1 | 0.151 | 1 | 0.151 |
| 2,2-Dichloropropane | U | <0.167 | <1.00 | <0.167 | µg/L | 1 | 0.167 | 1 | 0.167 |
| 1,2-Dichloroethane (EDC) | U | <0.0890 | <1.00 | <0.0890 | µg/L | 1 | 0.0890 | 1 | 0.089 |
| Chloroform | U | 1.07 | 1.07 | <0.121 | µg/L | 1 | 0.121 | 1 | 0.121 |
| 1,1,1-Trichloroethane | U | <0.170 | <1.00 | <0.170 | µg/L | 1 | 0.170 | 1 | 0.17 |
| 1,1-Dichloropropene | U | <0.156 | <1.00 | <0.156 | µg/L | 1 | 0.156 | 1 | 0.156 |
| Benzene | J | 0.288 | <1.00 | <0.113 | µg/L | 1 | 0.113 | 1 | 0.113 |
| Carbon Tetrachloride | U | <0.0930 | <1.00 | <0.0930 | µg/L | 1 | 0.0930 | 1 | 0.093 |
| 1,2-Dichloropropane | U | <0.129 | <1.00 | <0.129 | µg/L | 1 | 0.129 | 1 | 0.129 |
| Trichloroethene (TCE) | J | 77.8 | 77.8 | <0.160 | µg/L | 1 | 0.160 | 1 | 0.16 |
| Dibromomethane (methylene bromide) | U | <0.137 | <1.00 | <0.137 | µg/L | 1 | 0.137 | 1 | 0.137 |
| Bromodichloromethane | J | 0.165 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| 2-Chloroethyl vinyl ether | U R | <0.257 | <5.00 | <0.257 | µg/L | 1 | 0.257 | 5 | 0.257 |
| cis-1,3-Dichloropropene | U | <0.109 | <1.00 | <0.109 | µg/L | 1 | 0.109 | 1 | 0.109 |
| trans-1,3-Dichloropropene | U | <0.181 | <1.00 | <0.181 | µg/L | 1 | 0.181 | 1 | 0.181 |
| Toluene | U | <0.117 | <1.00 | <0.117 | µg/L | 1 | 0.117 | 1 | 0.117 |

continued ...

¹Concentration biased low.²Concentration biased low.

sample 209311 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|-----------------------------|------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| 1,1,2-Trichloroethane | U | <0.143 | <1.00 | <0.143 | µg/L | 1 | 0.143 | 1 | 0.143 |
| 1,3-Dichloropropane | U | <0.191 | <1.00 | <0.191 | µg/L | 1 | 0.191 | 1 | 0.191 |
| Dibromochloromethane | U | <0.118 | <1.00 | <0.118 | µg/L | 1 | 0.118 | 1 | 0.118 |
| 1,2-Dibromoethane (EDB) | U | <0.131 | <1.00 | <0.131 | µg/L | 1 | 0.131 | 1 | 0.131 |
| Tetrachloroethene (PCE) | U | <0.353 | <1.00 | <0.353 | µg/L | 1 | 0.353 | 1 | 0.353 |
| Chlorobenzene | U | <0.135 | <1.00 | <0.135 | µg/L | 1 | 0.135 | 1 | 0.135 |
| 1,1,1,2-Tetrachloroethane | U | <0.263 | <1.00 | <0.263 | µg/L | 1 | 0.263 | 1 | 0.263 |
| Ethylbenzene | U | <0.174 | <1.00 | <0.174 | µg/L | 1 | 0.174 | 1 | 0.174 |
| m,p-Xylene | U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Bromoform | U | <0.385 | <1.00 | <0.385 | µg/L | 1 | 0.385 | 1 | 0.385 |
| Styrene | U | <0.413 | <1.00 | <0.413 | µg/L | 1 | 0.413 | 1 | 0.413 |
| o-Xylene | U | <0.115 | <1.00 | <0.115 | µg/L | 1 | 0.115 | 1 | 0.115 |
| 1,1,2,2-Tetrachloroethane | U | <0.766 | <1.00 | <0.766 | µg/L | 1 | 0.766 | 1 | 0.766 |
| 2-Chlorotoluene | U | <0.132 | <1.00 | <0.132 | µg/L | 1 | 0.132 | 1 | 0.132 |
| 1,2,3-Trichloropropane | U | <0.599 | <1.00 | <0.599 | µg/L | 1 | 0.599 | 1 | 0.599 |
| Isopropylbenzene | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Bromobenzene | U | <0.266 | <1.00 | <0.266 | µg/L | 1 | 0.266 | 1 | 0.266 |
| n-Propylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,3,5-Trimethylbenzene | U | <0.124 | <1.00 | <0.124 | µg/L | 1 | 0.124 | 1 | 0.124 |
| tert-Butylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,2,4-Trimethylbenzene | U | <0.114 | <1.00 | <0.114 | µg/L | 1 | 0.114 | 1 | 0.114 |
| 1,4-Dichlorobenzene (para) | U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| sec-Butylbenzene | U | <0.452 | <1.00 | <0.452 | µg/L | 1 | 0.452 | 1 | 0.452 |
| 1,3-Dichlorobenzene (meta) | U | <0.214 | <1.00 | <0.214 | µg/L | 1 | 0.214 | 1 | 0.214 |
| p-Isopropyltoluene | U | <0.126 | <1.00 | <0.126 | µg/L | 1 | 0.126 | 1 | 0.126 |
| 4-Chlorotoluene | U | <0.161 | <1.00 | <0.161 | µg/L | 1 | 0.161 | 1 | 0.161 |
| 1,2-Dichlorobenzene (ortho) | U | <0.357 | <1.00 | <0.357 | µg/L | 1 | 0.357 | 1 | 0.357 |
| n-Butylbenzene | U | <0.125 | <1.00 | 0.163 | µg/L | 1 | 0.125 | 1 | 0.125 |
| 1,2-Dibromo-3-chloropropane | U | <0.977 | <5.00 | <0.977 | µg/L | 1 | 0.977 | 5 | 0.977 |
| 1,2,3-Trichlorobenzene | U | <0.400 | <5.00 | <0.400 | µg/L | 1 | 0.400 | 5 | 0.4 |
| 1,2,4-Trichlorobenzene | U | <0.227 | <5.00 | <0.227 | µg/L | 1 | 0.227 | 5 | 0.227 |
| Naphthalene | U | <0.672 | <5.00 | <0.672 | µg/L | 1 | 0.672 | 5 | 0.672 |
| Hexachlorobutadiene | U | <0.198 | <5.00 | 0.473 | µg/L | 1 | 0.198 | 5 | 0.198 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 49.8 | µg/L | 1 | 50.0 | 100 | 70 - 130 |
| Toluene-d8 | | 51.2 | µg/L | 1 | 50.0 | 102 | 70 - 130 |
| 4-Bromofluorobenzene (4-BFB) | | 51.2 | µg/L | 1 | 50.0 | 102 | 70 - 130 |

Sample: 209312 - HLSF-0154-HCF-001-0909

³Concentration biased low.⁴Concentration biased low.

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 63556

Prep Batch: 54254

Analytical Method: S 8260B

Date Analyzed: 2009-09-14

Sample Preparation: 2009-09-14

Prep Method: S 5030B

Analyzed By: JG

Prepared By: JG

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|----------------|--------------|--------------|--------------|-------|----------|--------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | U | <0.169 | <1.00 | <0.169 | µg/L | 1 | 0.169 | 1 | 0.169 |
| Dichlorodifluoromethane | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Chloromethane (methyl chloride) | U | <0.164 | <1.00 | <0.164 | µg/L | 1 | 0.164 | 1 | 0.164 |
| Vinyl Chloride | U | <0.110 | <1.00 | <0.110 | µg/L | 1 | 0.110 | 1 | 0.11 |
| Bromomethane (methyl bromide) | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |
| Chloroethane | U | <0.232 | <1.00 | <0.232 | µg/L | 1 | 0.232 | 1 | 0.232 |
| Trichlorofluoromethane | U | <0.149 | <1.00 | <0.149 | µg/L | 1 | 0.149 | 1 | 0.149 |
| Acetone | U ⁵ | <0.893 | <10.0 | 2.05 | µg/L | 1 | 0.893 | 10 | 0.893 |
| Iodomethane (methyl iodide) | U | <0.104 | <5.00 | <0.104 | µg/L | 1 | 0.104 | 5 | 0.104 |
| Carbon Disulfide | J | 0.530 | <1.00 | <0.120 | µg/L | 1 | 0.120 | 1 | 0.12 |
| Acrylonitrile | U | <0.189 | <1.00 | <0.189 | µg/L | 1 | 0.189 | 1 | 0.189 |
| 2-Butanone (MEK) | U ⁵ | <0.218 | <5.00 | 0.596 | µg/L | 1 | 0.218 | 5 | 0.218 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.554 | <5.00 | <0.554 | µg/L | 1 | 0.554 | 5 | 0.554 |
| 2-Hexanone | U | <0.863 | <5.00 | <0.863 | µg/L | 1 | 0.863 | 5 | 0.863 |
| trans-1,4-Dichloro-2-butene | U | <0.607 | <10.0 | <0.607 | µg/L | 1 | 0.607 | 10 | 0.607 |
| 1,1-Dichloroethene | U | <0.102 | <1.00 | <0.102 | µg/L | 1 | 0.102 | 1 | 0.102 |
| Methylene chloride | U | <0.270 | <5.00 | <0.270 | µg/L | 1 | 0.270 | 5 | 0.27 |
| MTBE | | 1.40 | 1.40 | <0.122 | µg/L | 1 | 0.122 | 1 | 0.122 |
| trans-1,2-Dichloroethene | U | <0.168 | <1.00 | <0.168 | µg/L | 1 | 0.168 | 1 | 0.168 |
| 1,1-Dichloroethane | J | 0.621 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| cis-1,2-Dichloroethene | U | <0.151 | <1.00 | <0.151 | µg/L | 1 | 0.151 | 1 | 0.151 |
| 2,2-Dichloropropane | U | <0.167 | <1.00 | <0.167 | µg/L | 1 | 0.167 | 1 | 0.167 |
| 1,2-Dichloroethane (EDC) | U | <0.0890 | <1.00 | <0.0890 | µg/L | 1 | 0.0890 | 1 | 0.089 |
| Chloroform | U | <0.121 | <1.00 | <0.121 | µg/L | 1 | 0.121 | 1 | 0.121 |
| 1,1,1-Trichloroethane | U | <0.170 | <1.00 | <0.170 | µg/L | 1 | 0.170 | 1 | 0.17 |
| 1,1-Dichloropropene | U | <0.156 | <1.00 | <0.156 | µg/L | 1 | 0.156 | 1 | 0.156 |
| Benzene | | 4.80 | 4.80 | <0.113 | µg/L | 1 | 0.113 | 1 | 0.113 |
| Carbon Tetrachloride | U | <0.0930 | <1.00 | <0.0930 | µg/L | 1 | 0.0930 | 1 | 0.093 |
| 1,2-Dichloropropane | U | <0.129 | <1.00 | <0.129 | µg/L | 1 | 0.129 | 1 | 0.129 |
| Trichloroethene (TCE) | J | 0.170 | <1.00 | <0.160 | µg/L | 1 | 0.160 | 1 | 0.16 |
| Dibromomethane (methylene bromide) | U | <0.137 | <1.00 | <0.137 | µg/L | 1 | 0.137 | 1 | 0.137 |
| Bromodichloromethane | U | <0.107 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| 2-Chloroethyl vinyl ether | U | <0.257 | <5.00 | <0.257 | µg/L | 1 | 0.257 | 5 | 0.257 |
| cis-1,3-Dichloropropene | U | <0.109 | <1.00 | <0.109 | µg/L | 1 | 0.109 | 1 | 0.109 |
| trans-1,3-Dichloropropene | U | <0.181 | <1.00 | <0.181 | µg/L | 1 | 0.181 | 1 | 0.181 |
| Toluene | J | 0.124 | <1.00 | <0.117 | µg/L | 1 | 0.117 | 1 | 0.117 |
| 1,1,2-Trichloroethane | U | <0.143 | <1.00 | <0.143 | µg/L | 1 | 0.143 | 1 | 0.143 |
| 1,3-Dichloropropane | U | <0.191 | <1.00 | <0.191 | µg/L | 1 | 0.191 | 1 | 0.191 |
| Dibromochloromethane | U | <0.118 | <1.00 | <0.118 | µg/L | 1 | 0.118 | 1 | 0.118 |
| 1,2-Dibromoethane (EDB) | U | <0.131 | <1.00 | <0.131 | µg/L | 1 | 0.131 | 1 | 0.131 |

*continued ...*⁵Concentration biased low.⁶Concentration biased low.

sample 209312 continued ...

| Parameter | Flag | SDL Based Result | SQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|-----------------------------|------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| Tetrachloroethene (PCE) | U | <0.353 | <1.00 | <0.353 | µg/L | 1 | 0.353 | 1 | 0.353 |
| Chlorobenzene | U | <0.135 | <1.00 | <0.135 | µg/L | 1 | 0.135 | 1 | 0.135 |
| 1,1,1,2-Tetrachloroethane | U | <0.263 | <1.00 | <0.263 | µg/L | 1 | 0.263 | 1 | 0.263 |
| Ethylbenzene | U | <0.174 | <1.00 | <0.174 | µg/L | 1 | 0.174 | 1 | 0.174 |
| m,p-Xylene | U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Bromoform | U | <0.385 | <1.00 | <0.385 | µg/L | 1 | 0.385 | 1 | 0.385 |
| Styrene | U | <0.413 | <1.00 | <0.413 | µg/L | 1 | 0.413 | 1 | 0.413 |
| o-Xylene | U | <0.115 | <1.00 | <0.115 | µg/L | 1 | 0.115 | 1 | 0.115 |
| 1,1,2,2-Tetrachloroethane | U | <0.766 | <1.00 | <0.766 | µg/L | 1 | 0.766 | 1 | 0.766 |
| 2-Chlorotoluene | U | <0.132 | <1.00 | <0.132 | µg/L | 1 | 0.132 | 1 | 0.132 |
| 1,2,3-Trichloropropane | U | <0.599 | <1.00 | <0.599 | µg/L | 1 | 0.599 | 1 | 0.599 |
| Isopropylbenzene | U | 2.22 | 2.22 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Bromobenzene | U | <0.266 | <1.00 | <0.266 | µg/L | 1 | 0.266 | 1 | 0.266 |
| n-Propylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,3,5-Trimethylbenzene | U | <0.124 | <1.00 | <0.124 | µg/L | 1 | 0.124 | 1 | 0.124 |
| tert-Butylbenzene | U | 0.195 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,2,4-Trimethylbenzene | U | <0.114 | <1.00 | <0.114 | µg/L | 1 | 0.114 | 1 | 0.114 |
| 1,4-Dichlorobenzene (para) | U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| sec-Butylbenzene | U | 3.21 | 3.21 | <0.452 | µg/L | 1 | 0.452 | 1 | 0.452 |
| 1,3-Dichlorobenzene (meta) | U | <0.214 | <1.00 | <0.214 | µg/L | 1 | 0.214 | 1 | 0.214 |
| p-Isopropyltoluene | U | <0.126 | <1.00 | <0.126 | µg/L | 1 | 0.126 | 1 | 0.126 |
| 4-Chlorotoluene | U | <0.161 | <1.00 | <0.161 | µg/L | 1 | 0.161 | 1 | 0.161 |
| 1,2-Dichlorobenzene (ortho) | U | <0.357 | <1.00 | <0.357 | µg/L | 1 | 0.357 | 1 | 0.357 |
| n-Butylbenzene | U | 0.696 | <1.00 | 0.163 | µg/L | 1 | 0.125 | 1 | 0.125 |
| 1,2-Dibromo-3-chloropropane | U | <0.977 | <5.00 | <0.977 | µg/L | 1 | 0.977 | 5 | 0.977 |
| 1,2,3-Trichlorobenzene | U | <0.400 | <5.00 | <0.400 | µg/L | 1 | 0.400 | 5 | 0.4 |
| 1,2,4-Trichlorobenzene | U | <0.227 | <5.00 | <0.227 | µg/L | 1 | 0.227 | 5 | 0.227 |
| Naphthalene | U | 35.0 | 35.0 | <0.672 | µg/L | 1 | 0.672 | 5 | 0.672 |
| Hexachlorobutadiene | U | <0.198 | <5.00 | 0.473 | µg/L | 1 | 0.198 | 5 | 0.198 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|-----------------|---------------------|--------------------|
| Dibromofluoromethane | | 48.6 | µg/L | 1 | 50.0 | 97 | 70 - 130 |
| Toluene-d8 | | 52.0 | µg/L | 1 | 50.0 | 104 | 70 - 130 |
| 4-Bromofluorobenzene (4-BFB) | | 51.1 | µg/L | 1 | 50.0 | 102 | 70 - 130 |

Sample: 209313 - HLSF-0154-DRW-005-0909

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 63556

Prep Batch: 54254

Analytical Method: S 8260B

Date Analyzed: 2009-09-14

Sample Preparation: 2009-09-14

Prep Method: S 5030B

Analyzed By: JG

Prepared By: JG

⁷Concentration biased low.⁸Concentration biased low.

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|-----------------------------|--------------|--------------|--------------|-------|----------|--------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | U | <0.169 | <1.00 | <0.169 | µg/L | 1 | 0.169 | 1 | 0.169 |
| Dichlorodifluoromethane | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Chloromethane (methyl chloride) | U | <0.164 | <1.00 | <0.164 | µg/L | 1 | 0.164 | 1 | 0.164 |
| Vinyl Chloride | U | <0.110 | <1.00 | <0.110 | µg/L | 1 | 0.110 | 1 | 0.11 |
| Bromomethane (methyl bromide) | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |
| Chloroethane | U | <0.232 | <1.00 | <0.232 | µg/L | 1 | 0.232 | 1 | 0.232 |
| Trichlorofluoromethane | U | <0.149 | <1.00 | <0.149 | µg/L | 1 | 0.149 | 1 | 0.149 |
| Acetone | J ⁹ U | 1.32 | <10.0 | 2.05 | µg/L | 1 | 0.893 | 10 | 0.893 |
| Iodomethane (methyl iodide) | U | <0.104 | <5.00 | <0.104 | µg/L | 1 | 0.104 | 5 | 0.104 |
| Carbon Disulfide | U | <0.120 | <1.00 | <0.120 | µg/L | 1 | 0.120 | 1 | 0.12 |
| Acrylonitrile | U | <0.189 | <1.00 | <0.189 | µg/L | 1 | 0.189 | 1 | 0.189 |
| 2-Butanone (MEK) | ¹⁰ U | <0.218 | <5.00 | 0.596 | µg/L | 1 | 0.218 | 5 | 0.218 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.554 | <5.00 | <0.554 | µg/L | 1 | 0.554 | 5 | 0.554 |
| 2-Hexanone | U | <0.863 | <5.00 | <0.863 | µg/L | 1 | 0.863 | 5 | 0.863 |
| trans 1,4-Dichloro-2-butene | U | <0.607 | <10.0 | <0.607 | µg/L | 1 | 0.607 | 10 | 0.607 |
| 1,1-Dichloroethene | J | 0.518 | <1.00 | <0.102 | µg/L | 1 | 0.102 | 1 | 0.102 |
| Methylene chloride | U | <0.270 | <5.00 | <0.270 | µg/L | 1 | 0.270 | 5 | 0.27 |
| MTBE | | 2.42 | 2.42 | <0.122 | µg/L | 1 | 0.122 | 1 | 0.122 |
| trans-1,2-Dichloroethene | U | <0.168 | <1.00 | <0.168 | µg/L | 1 | 0.168 | 1 | 0.168 |
| 1,1-Dichloroethane | | 42.0 | 42.0 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| cis-1,2-Dichloroethene | U | <0.151 | <1.00 | <0.151 | µg/L | 1 | 0.151 | 1 | 0.151 |
| 2,2-Dichloropropane | U | <0.167 | <1.00 | <0.167 | µg/L | 1 | 0.167 | 1 | 0.167 |
| 1,2-Dichloroethane (EDC) | U | <0.0890 | <1.00 | <0.0890 | µg/L | 1 | 0.0890 | 1 | 0.089 |
| Chloroform | U | <0.121 | <1.00 | <0.121 | µg/L | 1 | 0.121 | 1 | 0.121 |
| 1,1,1-Trichloroethane | U | <0.170 | <1.00 | <0.170 | µg/L | 1 | 0.170 | 1 | 0.17 |
| 1,1-Dichloropropene | U | <0.156 | <1.00 | <0.156 | µg/L | 1 | 0.156 | 1 | 0.156 |
| Benzene | | 5.49 | 5.49 | <0.113 | µg/L | 1 | 0.113 | 1 | 0.113 |
| Carbon Tetrachloride | U | <0.0930 | <1.00 | <0.0930 | µg/L | 1 | 0.0930 | 1 | 0.093 |
| 1,2-Dichloropropane | U | <0.129 | <1.00 | <0.129 | µg/L | 1 | 0.129 | 1 | 0.129 |
| Trichloroethene (TCE) | J | 0.529 | <1.00 | <0.160 | µg/L | 1 | 0.160 | 1 | 0.16 |
| Dibromomethane (methylene bromide) | U | <0.137 | <1.00 | <0.137 | µg/L | 1 | 0.137 | 1 | 0.137 |
| Bromodichloromethane | U | <0.107 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| 2-Chloroethyl vinyl ether | U | <0.257 | <5.00 | <0.257 | µg/L | 1 | 0.257 | 5 | 0.257 |
| cis-1,3-Dichloropropene | U | <0.109 | <1.00 | <0.109 | µg/L | 1 | 0.109 | 1 | 0.109 |
| trans-1,3-Dichloropropene | U | <0.181 | <1.00 | <0.181 | µg/L | 1 | 0.181 | 1 | 0.181 |
| Toluene | J | 0.231 | <1.00 | <0.117 | µg/L | 1 | 0.117 | 1 | 0.117 |
| 1,1,2-Trichloroethane | U | <0.143 | <1.00 | <0.143 | µg/L | 1 | 0.143 | 1 | 0.143 |
| 1,3-Dichloropropane | U | <0.191 | <1.00 | <0.191 | µg/L | 1 | 0.191 | 1 | 0.191 |
| Dibromochloromethane | U | <0.118 | <1.00 | <0.118 | µg/L | 1 | 0.118 | 1 | 0.118 |
| 1,2-Dibromoethane (EDB) | U | <0.131 | <1.00 | <0.131 | µg/L | 1 | 0.131 | 1 | 0.131 |
| Tetrachloroethene (PCE) | ¹¹ U J | <0.353 | <1.00 | <0.353 | µg/L | 1 | 0.353 | 1 | 0.353 |
| Chlorobenzene | U | <0.135 | <1.00 | <0.135 | µg/L | 1 | 0.135 | 1 | 0.135 |
| 1,1,1,2-Tetrachloroethane | U | <0.263 | <1.00 | <0.263 | µg/L | 1 | 0.263 | 1 | 0.263 |
| Ethylbenzene | U | <0.174 | <1.00 | <0.174 | µg/L | 1 | 0.174 | 1 | 0.174 |

continued ...

⁹Concentration biased low.¹⁰Concentration biased low.¹¹Concentration biased low.

sample 209313 continued ...

| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|-----------------------------|----------------------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| m,p-Xylene | U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Bromoform | U | <0.385 | <1.00 | <0.385 | µg/L | 1 | 0.385 | 1 | 0.385 |
| Styrene | U | <0.413 | <1.00 | <0.413 | µg/L | 1 | 0.413 | 1 | 0.413 |
| o-Xylene | J | 0.179 | <1.00 | <0.115 | µg/L | 1 | 0.115 | 1 | 0.115 |
| 1,1,2,2-Tetrachloroethane | U | <0.766 | <1.00 | <0.766 | µg/L | 1 | 0.766 | 1 | 0.766 |
| 2-Chlorotoluene | U | <0.132 | <1.00 | <0.132 | µg/L | 1 | 0.132 | 1 | 0.132 |
| 1,2,3-Trichloropropane | U | <0.599 | <1.00 | <0.599 | µg/L | 1 | 0.599 | 1 | 0.599 |
| Isopropylbenzene | | 1.49 | 1.49 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Bromobenzene | U | <0.266 | <1.00 | <0.266 | µg/L | 1 | 0.266 | 1 | 0.266 |
| n-Propylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,3,5-Trimethylbenzene | U | <0.124 | <1.00 | <0.124 | µg/L | 1 | 0.124 | 1 | 0.124 |
| tert-Butylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,2,4-Trimethylbenzene | U ^{UB} | 0.123 | <1.00 | <0.114 | µg/L | 1 | 0.114 | 1 | 0.114 |
| 1,4-Dichlorobenzene (para) | U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| sec-Butylbenzene | | 2.20 | 2.20 | <0.452 | µg/L | 1 | 0.452 | 1 | 0.452 |
| 1,3-Dichlorobenzene (meta) | U | <0.214 | <1.00 | <0.214 | µg/L | 1 | 0.214 | 1 | 0.214 |
| p-Isopropyltoluene | U | <0.126 | <1.00 | <0.126 | µg/L | 1 | 0.126 | 1 | 0.126 |
| 4-Chlorotoluene | U | <0.161 | <1.00 | <0.161 | µg/L | 1 | 0.161 | 1 | 0.161 |
| 1,2-Dichlorobenzene (ortho) | U | <0.357 | <1.00 | <0.357 | µg/L | 1 | 0.357 | 1 | 0.357 |
| n-Butylbenzene | JB | 0.631 | <1.00 | 0.163 | µg/L | 1 | 0.125 | 1 | 0.125 |
| 1,2-Dibromo-3-chloropropane | U | <0.977 | <5.00 | <0.977 | µg/L | 1 | 0.977 | 5 | 0.977 |
| 1,2,3-Trichlorobenzene | U | <0.400 | <5.00 | <0.400 | µg/L | 1 | 0.400 | 5 | 0.4 |
| 1,2,4-Trichlorobenzene | U | <0.227 | <5.00 | <0.227 | µg/L | 1 | 0.227 | 5 | 0.227 |
| Naphthalene | | 9.47 | 9.47 | <0.672 | µg/L | 1 | 0.672 | 5 | 0.672 |
| Hexachlorobutadiene | U | <0.198 | <5.00 | 0.473 | µg/L | 1 | 0.198 | 5 | 0.198 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|-----------------|---------------------|--------------------|
| Dibromofluoromethane | | 48.9 | µg/L | 1 | 50.0 | 98 | 70 - 130 |
| Toluene-d8 | | 50.8 | µg/L | 1 | 50.0 | 102 | 70 - 130 |
| 4-Bromofluorobenzene (4-BFB) | | 51.4 | µg/L | 1 | 50.0 | 103 | 70 - 130 |

Sample: 209314 - HLSF-0143-HMW-013-0909

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 63556

Prep Batch: 54254

Analytical Method: S 8260B

Date Analyzed: 2009-09-14

Sample Preparation: 2009-09-14

Prep Method: S 5030B

Analyzed By: JG

Prepared By: JG

continued ...

¹²Concentration biased low.

sample 209314 continued ...

| Parameter | Flag | SDL Based Result | SQL Based Result | Method Blank Result | Units | Dilution | SDL | SQL (Unadjusted) | MDL (Unadjusted) |
|------------------------------------|------|------------------------|------------------------|---------------------------|-------|----------|--------|---------------------|---------------------|
| Parameter | Flag | SDL Based Result | SQL Based Result | Method Blank Result | Units | Dilution | SDL | SQL (Unadjusted) | MDL (Unadjusted) |
| Bromochloromethane | U | <0.169 | <1.00 | <0.169 | µg/L | 1 | 0.169 | 1 | 0.169 |
| Dichlorodifluoromethane | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Chloromethane (methyl chloride) | U | <0.164 | <1.00 | <0.164 | µg/L | 1 | 0.164 | 1 | 0.164 |
| Vinyl Chloride | U | <0.110 | <1.00 | <0.110 | µg/L | 1 | 0.110 | 1 | 0.11 |
| Bromomethane (methyl bromide) | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |
| Chloroethane | U | <0.232 | <1.00 | <0.232 | µg/L | 1 | 0.232 | 1 | 0.232 |
| Trichlorofluoromethane | U | <0.149 | <1.00 | <0.149 | µg/L | 1 | 0.149 | 1 | 0.149 |
| Acetone | U | <0.893 | <10.0 | 2.05 | µg/L | 1 | 0.893 | 10 | 0.893 |
| Iodomethane (methyl iodide) | U | <0.104 | <5.00 | <0.104 | µg/L | 1 | 0.104 | 5 | 0.104 |
| Carbon Disulfide | U | <0.120 | <1.00 | <0.120 | µg/L | 1 | 0.120 | 1 | 0.12 |
| Acrylonitrile | U | <0.189 | <1.00 | <0.189 | µg/L | 1 | 0.189 | 1 | 0.189 |
| 2-Butanone (MEK) | U | <0.218 | <5.00 | 0.596 | µg/L | 1 | 0.218 | 5 | 0.218 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.554 | <5.00 | <0.554 | µg/L | 1 | 0.554 | 5 | 0.554 |
| 2-Hexanone | U | <0.863 | <5.00 | <0.863 | µg/L | 1 | 0.863 | 5 | 0.863 |
| trans 1,4-Dichloro-2-butene | U | <0.607 | <10.0 | <0.607 | µg/L | 1 | 0.607 | 10 | 0.607 |
| 1,1-Dichloroethene | U | <0.102 | <1.00 | <0.102 | µg/L | 1 | 0.102 | 1 | 0.102 |
| Methylene chloride | U | <0.270 | <5.00 | <0.270 | µg/L | 1 | 0.270 | 5 | 0.27 |
| MTBE | U | <0.122 | <1.00 | <0.122 | µg/L | 1 | 0.122 | 1 | 0.122 |
| trans-1,2-Dichloroethene | U | <0.168 | <1.00 | <0.168 | µg/L | 1 | 0.168 | 1 | 0.168 |
| 1,1-Dichloroethane | U | 1.16 | 1.16 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| cis-1,2-Dichloroethene | U | <0.151 | <1.00 | <0.151 | µg/L | 1 | 0.151 | 1 | 0.151 |
| 2,2-Dichloropropane | U | <0.167 | <1.00 | <0.167 | µg/L | 1 | 0.167 | 1 | 0.167 |
| 1,2-Dichloroethane (EDC) | U | <0.0890 | <1.00 | <0.0890 | µg/L | 1 | 0.0890 | 1 | 0.089 |
| Chloroform | U | 0.982 | <1.00 | <0.121 | µg/L | 1 | 0.121 | 1 | 0.121 |
| 1,1,1-Trichloroethane | U | <0.170 | <1.00 | <0.170 | µg/L | 1 | 0.170 | 1 | 0.17 |
| 1,1-Dichloropropene | U | <0.156 | <1.00 | <0.156 | µg/L | 1 | 0.156 | 1 | 0.156 |
| Benzene | U | <0.113 | <1.00 | <0.113 | µg/L | 1 | 0.113 | 1 | 0.113 |
| Carbon Tetrachloride | U | <0.0930 | <1.00 | <0.0930 | µg/L | 1 | 0.0930 | 1 | 0.093 |
| 1,2-Dichloropropane | U | <0.129 | <1.00 | <0.129 | µg/L | 1 | 0.129 | 1 | 0.129 |
| Trichloroethene (TCE) | U | <0.160 | <1.00 | <0.160 | µg/L | 1 | 0.160 | 1 | 0.16 |
| Dibromomethane (methylene bromide) | U | <0.137 | <1.00 | <0.137 | µg/L | 1 | 0.137 | 1 | 0.137 |
| Bromodichloromethane | U | <0.107 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| 2-Chloroethyl vinyl ether | U | <0.257 | <5.00 | <0.257 | µg/L | 1 | 0.257 | 5 | 0.257 |
| cis-1,3-Dichloropropene | U | <0.109 | <1.00 | <0.109 | µg/L | 1 | 0.109 | 1 | 0.109 |
| trans-1,3-Dichloropropene | U | <0.181 | <1.00 | <0.181 | µg/L | 1 | 0.181 | 1 | 0.181 |
| Toluene | U | <0.117 | <1.00 | <0.117 | µg/L | 1 | 0.117 | 1 | 0.117 |
| 1,1,2-Trichloroethane | U | <0.143 | <1.00 | <0.143 | µg/L | 1 | 0.143 | 1 | 0.143 |
| 1,3-Dichloropropane | U | <0.191 | <1.00 | <0.191 | µg/L | 1 | 0.191 | 1 | 0.191 |
| Dibromochloromethane | U | <0.118 | <1.00 | <0.118 | µg/L | 1 | 0.118 | 1 | 0.118 |

continued ...

¹³Concentration biased low.¹⁴Concentration biased low.

sample 209314 continued ...

| Parameter | Flag | SDL Based Result | SQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|-----------------------------|------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| 1,2-Dibromoethane (EDB) | U | <0.131 | <1.00 | <0.131 | µg/L | 1 | 0.131 | 1 | 0.131 |
| Tetrachloroethene (PCE) | U | <0.353 | <1.00 | <0.353 | µg/L | 1 | 0.353 | 1 | 0.353 |
| Chlorobenzene | U | <0.135 | <1.00 | <0.135 | µg/L | 1 | 0.135 | 1 | 0.135 |
| 1,1,1,2-Tetrachloroethane | U | <0.263 | <1.00 | <0.263 | µg/L | 1 | 0.263 | 1 | 0.263 |
| Ethylbenzene | U | <0.174 | <1.00 | <0.174 | µg/L | 1 | 0.174 | 1 | 0.174 |
| m,p-Xylene | U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Bromoform | U | <0.385 | <1.00 | <0.385 | µg/L | 1 | 0.385 | 1 | 0.385 |
| Styrene | U | <0.413 | <1.00 | <0.413 | µg/L | 1 | 0.413 | 1 | 0.413 |
| o-Xylene | U | <0.115 | <1.00 | <0.115 | µg/L | 1 | 0.115 | 1 | 0.115 |
| 1,1,2,2-Tetrachloroethane | U | <0.766 | <1.00 | <0.766 | µg/L | 1 | 0.766 | 1 | 0.766 |
| 2-Chlorotoluene | U | <0.132 | <1.00 | <0.132 | µg/L | 1 | 0.132 | 1 | 0.132 |
| 1,2,3-Trichloropropane | U | <0.599 | <1.00 | <0.599 | µg/L | 1 | 0.599 | 1 | 0.599 |
| Isopropylbenzene | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Bromobenzene | U | <0.266 | <1.00 | <0.266 | µg/L | 1 | 0.266 | 1 | 0.266 |
| n-Propylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,3,5-Trimethylbenzene | U | <0.124 | <1.00 | <0.124 | µg/L | 1 | 0.124 | 1 | 0.124 |
| tert-Butylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,2,4-Trimethylbenzene | U | <0.114 | <1.00 | <0.114 | µg/L | 1 | 0.114 | 1 | 0.114 |
| 1,4-Dichlorobenzene (para) | U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| sec-Butylbenzene | U | <0.452 | <1.00 | <0.452 | µg/L | 1 | 0.452 | 1 | 0.452 |
| 1,3-Dichlorobenzene (meta) | U | <0.214 | <1.00 | <0.214 | µg/L | 1 | 0.214 | 1 | 0.214 |
| p-Isopropyltoluene | U | <0.126 | <1.00 | <0.126 | µg/L | 1 | 0.126 | 1 | 0.126 |
| 4-Chlorotoluene | U | <0.161 | <1.00 | <0.161 | µg/L | 1 | 0.161 | 1 | 0.161 |
| 1,2-Dichlorobenzene (ortho) | U | <0.357 | <1.00 | <0.357 | µg/L | 1 | 0.357 | 1 | 0.357 |
| n-Butylbenzene | U | <0.125 | <1.00 | 0.163 | µg/L | 1 | 0.125 | 1 | 0.125 |
| 1,2-Dibromo-3-chloropropane | U | <0.977 | <5.00 | <0.977 | µg/L | 1 | 0.977 | 5 | 0.977 |
| 1,2,3-Trichlorobenzene | U | <0.400 | <5.00 | <0.400 | µg/L | 1 | 0.400 | 5 | 0.4 |
| 1,2,4-Trichlorobenzene | U | <0.227 | <5.00 | <0.227 | µg/L | 1 | 0.227 | 5 | 0.227 |
| Naphthalene | U | <0.672 | <5.00 | <0.672 | µg/L | 1 | 0.672 | 5 | 0.672 |
| Hexachlorobutadiene | U | <0.198 | <5.00 | 0.473 | µg/L | 1 | 0.198 | 5 | 0.198 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|-----------------|---------------------|--------------------|
| Dibromofluoromethane | | 48.8 | µg/L | 1 | 50.0 | 98 | 70 - 130 |
| Toluene-d8 | | 51.6 | µg/L | 1 | 50.0 | 103 | 70 - 130 |
| 4-Bromofluorobenzene (4-BFB) | | 51.0 | µg/L | 1 | 50.0 | 102 | 70 - 130 |

Sample: 209315 - HLSF-0143-HMW-036-0909

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 63556

Analytical Method: S 8260B

Date Analyzed: 2009-09-14

Prep Method: S 5030B

Analyzed By: JG

¹⁵Concentration biased low.¹⁶Concentration biased low.

Prep Batch: 54254

Sample Preparation: 2009-09-14

Prepared By: JG

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|------------------------------------|------|--------------|--------------|--------------|-------|----------|--------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| Bromochloromethane | U | <0.169 | <1.00 | <0.169 | µg/L | 1 | 0.169 | 1 | 0.169 |
| Dichlorodifluoromethane | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Chloromethane (methyl chloride) | U | <0.164 | <1.00 | <0.164 | µg/L | 1 | 0.164 | 1 | 0.164 |
| Vinyl Chloride | U | <0.110 | <1.00 | <0.110 | µg/L | 1 | 0.110 | 1 | 0.11 |
| Bromomethane (methyl bromide) | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |
| Chloroethane | U | <0.232 | <1.00 | <0.232 | µg/L | 1 | 0.232 | 1 | 0.232 |
| Trichlorofluoromethane | U | <0.149 | <1.00 | <0.149 | µg/L | 1 | 0.149 | 1 | 0.149 |
| Acetone | U | 1.15 | <10.0 | 2.05 | µg/L | 1 | 0.893 | 10 | 0.893 |
| Iodomethane (methyl iodide) | U | <0.104 | <5.00 | <0.104 | µg/L | 1 | 0.104 | 5 | 0.104 |
| Carbon Disulfide | U | <0.120 | <1.00 | <0.120 | µg/L | 1 | 0.120 | 1 | 0.12 |
| Acrylonitrile | U | <0.189 | <1.00 | <0.189 | µg/L | 1 | 0.189 | 1 | 0.189 |
| 2-Butanone (MEK) | U | <0.218 | <5.00 | 0.596 | µg/L | 1 | 0.218 | 5 | 0.218 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.554 | <5.00 | <0.554 | µg/L | 1 | 0.554 | 5 | 0.554 |
| 2-Hexanone | U | <0.863 | <5.00 | <0.863 | µg/L | 1 | 0.863 | 5 | 0.863 |
| trans-1,4-Dichloro-2-butene | U | <0.607 | <10.0 | <0.607 | µg/L | 1 | 0.607 | 10 | 0.607 |
| 1,1-Dichloroethene | U | <0.102 | <1.00 | <0.102 | µg/L | 1 | 0.102 | 1 | 0.102 |
| Methylene chloride | J | 0.283 | <5.00 | <0.270 | µg/L | 1 | 0.270 | 5 | 0.27 |
| MTBE | U | <0.122 | <1.00 | <0.122 | µg/L | 1 | 0.122 | 1 | 0.122 |
| trans-1,2-Dichloroethene | U | <0.168 | <1.00 | <0.168 | µg/L | 1 | 0.168 | 1 | 0.168 |
| 1,1-Dichloroethane | J | 0.273 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| cis-1,2-Dichloroethene | U | <0.151 | <1.00 | <0.151 | µg/L | 1 | 0.151 | 1 | 0.151 |
| 2,2-Dichloropropane | U | <0.167 | <1.00 | <0.167 | µg/L | 1 | 0.167 | 1 | 0.167 |
| 1,2-Dichloroethane (EDC) | U | <0.0890 | <1.00 | <0.0890 | µg/L | 1 | 0.0890 | 1 | 0.089 |
| Chloroform | J | 0.549 | <1.00 | <0.121 | µg/L | 1 | 0.121 | 1 | 0.121 |
| 1,1,1-Trichloroethane | U | <0.170 | <1.00 | <0.170 | µg/L | 1 | 0.170 | 1 | 0.17 |
| 1,1-Dichloropropene | U | <0.156 | <1.00 | <0.156 | µg/L | 1 | 0.156 | 1 | 0.156 |
| Benzene | J | 0.130 | <1.00 | <0.113 | µg/L | 1 | 0.113 | 1 | 0.113 |
| Carbon Tetrachloride | U | <0.0930 | <1.00 | <0.0930 | µg/L | 1 | 0.0930 | 1 | 0.093 |
| 1,2-Dichloropropane | U | <0.129 | <1.00 | <0.129 | µg/L | 1 | 0.129 | 1 | 0.129 |
| Trichloroethene (TCE) | U | <0.160 | <1.00 | <0.160 | µg/L | 1 | 0.160 | 1 | 0.16 |
| Dibromomethane (methylene bromide) | U | <0.137 | <1.00 | <0.137 | µg/L | 1 | 0.137 | 1 | 0.137 |
| Bromodichloromethane | U | <0.107 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| 2-Chloroethyl vinyl ether | U | <0.257 | <5.00 | <0.257 | µg/L | 1 | 0.257 | 5 | 0.257 |
| cis-1,3-Dichloropropene | U | <0.109 | <1.00 | <0.109 | µg/L | 1 | 0.109 | 1 | 0.109 |
| trans-1,3-Dichloropropene | U | <0.181 | <1.00 | <0.181 | µg/L | 1 | 0.181 | 1 | 0.181 |
| Toluene | U | <0.117 | <1.00 | <0.117 | µg/L | 1 | 0.117 | 1 | 0.117 |
| 1,1,2-Trichloroethane | U | <0.143 | <1.00 | <0.143 | µg/L | 1 | 0.143 | 1 | 0.143 |
| 1,3-Dichloropropane | U | <0.191 | <1.00 | <0.191 | µg/L | 1 | 0.191 | 1 | 0.191 |
| Dibromochloromethane | U | <0.118 | <1.00 | <0.118 | µg/L | 1 | 0.118 | 1 | 0.118 |
| 1,2-Dibromoethane (EDB) | U | <0.131 | <1.00 | <0.131 | µg/L | 1 | 0.131 | 1 | 0.131 |
| Tetrachloroethene (PCE) | U | <0.353 | <1.00 | <0.353 | µg/L | 1 | 0.353 | 1 | 0.353 |
| Chlorobenzene | U | <0.135 | <1.00 | <0.135 | µg/L | 1 | 0.135 | 1 | 0.135 |

continued ...

¹⁷Concentration biased low.¹⁸Concentration biased low.¹⁹Concentration biased low.

sample 209315 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|-----------------------------|------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| 1,1,1,2-Tetrachloroethane | U | <0.263 | <1.00 | <0.263 | µg/L | 1 | 0.263 | 1 | 0.263 |
| Ethylbenzene | U | <0.174 | <1.00 | <0.174 | µg/L | 1 | 0.174 | 1 | 0.174 |
| m,p-Xylene | U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Bromoform | U | <0.385 | <1.00 | <0.385 | µg/L | 1 | 0.385 | 1 | 0.385 |
| Styrene | U | <0.413 | <1.00 | <0.413 | µg/L | 1 | 0.413 | 1 | 0.413 |
| o-Xylene | J | 0.122 | <1.00 | <0.115 | µg/L | 1 | 0.115 | 1 | 0.115 |
| 1,1,2,2-Tetrachloroethane | U | <0.766 | <1.00 | <0.766 | µg/L | 1 | 0.766 | 1 | 0.766 |
| 2-Chlorotoluene | U | <0.132 | <1.00 | <0.132 | µg/L | 1 | 0.132 | 1 | 0.132 |
| 1,2,3-Trichloropropane | U | <0.599 | <1.00 | <0.599 | µg/L | 1 | 0.599 | 1 | 0.599 |
| Isopropylbenzene | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Bromobenzene | U | <0.266 | <1.00 | <0.266 | µg/L | 1 | 0.266 | 1 | 0.266 |
| n-Propylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,3,5-Trimethylbenzene | U | <0.124 | <1.00 | <0.124 | µg/L | 1 | 0.124 | 1 | 0.124 |
| tert-Butylbenzene | U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,2,4-Trimethylbenzene | 20 U | <0.114 | <1.00 | <0.114 | µg/L | 1 | 0.114 | 1 | 0.114 |
| 1,4-Dichlorobenzene (para) | U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| sec-Butylbenzene | U | <0.452 | <1.00 | <0.452 | µg/L | 1 | 0.452 | 1 | 0.452 |
| 1,3-Dichlorobenzene (meta) | U | <0.214 | <1.00 | <0.214 | µg/L | 1 | 0.214 | 1 | 0.214 |
| p-Isopropyltoluene | U | <0.126 | <1.00 | <0.126 | µg/L | 1 | 0.126 | 1 | 0.126 |
| 4-Chlorotoluene | U | <0.161 | <1.00 | <0.161 | µg/L | 1 | 0.161 | 1 | 0.161 |
| 1,2-Dichlorobenzene (ortho) | U | <0.357 | <1.00 | <0.357 | µg/L | 1 | 0.357 | 1 | 0.357 |
| n-Butylbenzene | U | <0.125 | <1.00 | 0.163 | µg/L | 1 | 0.125 | 1 | 0.125 |
| 1,2-Dibromo-3-chloropropane | U | <0.977 | <5.00 | <0.977 | µg/L | 1 | 0.977 | 5 | 0.977 |
| 1,2,3-Trichlorobenzene | U | <0.400 | <5.00 | <0.400 | µg/L | 1 | 0.400 | 5 | 0.4 |
| 1,2,4-Trichlorobenzene | U | <0.227 | <5.00 | <0.227 | µg/L | 1 | 0.227 | 5 | 0.227 |
| Naphthalene | U | <0.672 | <5.00 | <0.672 | µg/L | 1 | 0.672 | 5 | 0.672 |
| Hexachlorobutadiene | U | <0.198 | <5.00 | 0.473 | µg/L | 1 | 0.198 | 5 | 0.198 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 50.4 | µg/L | 1 | 50.0 | 101 | 70 - 130 |
| Toluene-d8 | | 51.6 | µg/L | 1 | 50.0 | 103 | 70 - 130 |
| 4-Bromofluorobenzene (4-BFB) | | 51.8 | µg/L | 1 | 50.0 | 104 | 70 - 130 |

Sample: 209316 - HLSF-0154-DRW-004-0909

Laboratory: Lubbock,

Analysis: Volatiles

QC Batch: 63556

Prep Batch: 54254

Analytical Method: S 8260B

Date Analyzed: 2009-09-14

Sample Preparation: 2009-09-14

Prep Method: S 5030B

Analyzed By: JG

Prepared By: JG

²⁰Concentration biased low.

| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|------------------------------------|-----------------|------------------------|------------------------|---------------------------|-------|----------|-------|---------------------|---------------------|
| Bromochloromethane | U | <0.845 | <5.00 | <0.845 | µg/L | 5 | 0.845 | 1 | 0.169 |
| Dichlorodifluoromethane | U | <0.725 | <5.00 | <0.725 | µg/L | 5 | 0.725 | 1 | 0.145 |
| Chloromethane (methyl chloride) | U | <0.820 | <5.00 | <0.820 | µg/L | 5 | 0.820 | 1 | 0.164 |
| Vinyl Chloride | U | <0.550 | <5.00 | <0.550 | µg/L | 5 | 0.550 | 1 | 0.11 |
| Bromomethane (methyl bromide) | U | <2.70 | <25.0 | <2.70 | µg/L | 5 | 2.70 | 5 | 0.54 |
| Chloroethane | U | <1.16 | <5.00 | <1.16 | µg/L | 5 | 1.16 | 1 | 0.232 |
| Trichlorofluoromethane | U | <0.745 | <5.00 | <0.745 | µg/L | 5 | 0.745 | 1 | 0.149 |
| Acetone | J ²¹ | 8.02 | <50.0 | 10.2 | µg/L | 5 | 4.46 | 10 | 0.893 |
| Iodomethane (methyl iodide) | U | <0.520 | <25.0 | <0.520 | µg/L | 5 | 0.520 | 5 | 0.104 |
| Carbon Disulfide | U | <0.600 | <5.00 | <0.600 | µg/L | 5 | 0.600 | 1 | 0.12 |
| Acrylonitrile | U | <0.945 | <5.00 | <0.945 | µg/L | 5 | 0.945 | 1 | 0.189 |
| 2-Butanone (MEK) | J ²² | 2.24 | <25.0 | 2.98 | µg/L | 5 | 1.09 | 5 | 0.218 |
| 4-Methyl-2-pentanone (MIBK) | U | <2.77 | <25.0 | <2.77 | µg/L | 5 | 2.77 | 5 | 0.554 |
| 2-Hexanone | U | <4.32 | <25.0 | <4.32 | µg/L | 5 | 4.32 | 5 | 0.863 |
| trans-1,4-Dichloro-2-butene | U | <3.04 | <50.0 | <3.04 | µg/L | 5 | 3.04 | 10 | 0.607 |
| 1,1-Dichloroethene | J | 4.01 | <5.00 | <0.510 | µg/L | 5 | 0.510 | 1 | 0.102 |
| Methylene chloride | J | 1.45 | <25.0 | <1.35 | µg/L | 5 | 1.35 | 5 | 0.27 |
| MTBE | | 7.13 | 7.13 | <0.610 | µg/L | 5 | 0.610 | 1 | 0.122 |
| trans-1,2-Dichloroethene | U | <0.840 | <5.00 | <0.840 | µg/L | 5 | 0.840 | 1 | 0.168 |
| 1,1-Dichloroethane | | 142 | 142 | <0.535 | µg/L | 5 | 0.535 | 1 | 0.107 |
| cis-1,2-Dichloroethene | U | <0.755 | <5.00 | <0.755 | µg/L | 5 | 0.755 | 1 | 0.151 |
| 2,2-Dichloropropane | U | <0.835 | <5.00 | <0.835 | µg/L | 5 | 0.835 | 1 | 0.167 |
| 1,2-Dichloroethane (EDC) | U | <0.445 | <5.00 | <0.445 | µg/L | 5 | 0.445 | 1 | 0.089 |
| Chloroform | U | <0.605 | <5.00 | <0.605 | µg/L | 5 | 0.605 | 1 | 0.121 |
| 1,1,1-Trichloroethane | U | <0.850 | <5.00 | <0.850 | µg/L | 5 | 0.850 | 1 | 0.17 |
| 1,1-Dichloropropene | | 7.60 | 7.60 | <0.780 | µg/L | 5 | 0.780 | 1 | 0.156 |
| Benzene | | 5.10 | 5.10 | <0.565 | µg/L | 5 | 0.565 | 1 | 0.113 |
| Carbon Tetrachloride | U | <0.465 | <5.00 | <0.465 | µg/L | 5 | 0.465 | 1 | 0.093 |
| 1,2-Dichloropropane | U | <0.645 | <5.00 | <0.645 | µg/L | 5 | 0.645 | 1 | 0.129 |
| Trichloroethene (TCE) | U | <0.800 | <5.00 | <0.800 | µg/L | 5 | 0.800 | 1 | 0.16 |
| Dibromomethane (methylene bromide) | U | <0.685 | <5.00 | <0.685 | µg/L | 5 | 0.685 | 1 | 0.137 |
| Bromodichloromethane | U | <0.535 | <5.00 | <0.535 | µg/L | 5 | 0.535 | 1 | 0.107 |
| 2-Chloroethyl vinyl ether | U | <1.28 | <25.0 | <1.28 | µg/L | 5 | 1.28 | 5 | 0.257 |
| cis-1,3-Dichloropropene | U | <0.545 | <5.00 | <0.545 | µg/L | 5 | 0.545 | 1 | 0.109 |
| trans-1,3-Dichloropropene | U | <0.905 | <5.00 | <0.905 | µg/L | 5 | 0.905 | 1 | 0.181 |
| Toluene | J | 0.607 | <5.00 | <0.585 | µg/L | 5 | 0.585 | 1 | 0.117 |
| 1,1,2-Trichloroethane | U | <0.715 | <5.00 | <0.715 | µg/L | 5 | 0.715 | 1 | 0.143 |
| 1,3-Dichloropropane | U | <0.955 | <5.00 | <0.955 | µg/L | 5 | 0.955 | 1 | 0.191 |
| Dibromochloromethane | U | <0.590 | <5.00 | <0.590 | µg/L | 5 | 0.590 | 1 | 0.118 |
| 1,2-Dibromoethane (EDB) | U | <0.655 | <5.00 | <0.655 | µg/L | 5 | 0.655 | 1 | 0.131 |
| Tetrachloroethene (PCE) | J ²³ | <1.76 | <5.00 | <1.76 | µg/L | 5 | 1.76 | 1 | 0.353 |
| Chlorobenzene | U | <0.675 | <5.00 | <0.675 | µg/L | 5 | 0.675 | 1 | 0.135 |
| 1,1,1,2-Tetrachloroethane | U | <1.32 | <5.00 | <1.32 | µg/L | 5 | 1.32 | 1 | 0.263 |
| Ethylbenzene | U | <0.870 | <5.00 | <0.870 | µg/L | 5 | 0.870 | 1 | 0.174 |

continued ...

²¹Concentration biased low.²²Concentration biased low.²³Concentration biased low.

sample 209316 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|-----------------------------|------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| m,p-Xylene | U | <1.10 | <5.00 | <1.10 | µg/L | 5 | 1.10 | 1 | 0.22 |
| Bromoform | U | <1.92 | <5.00 | <1.92 | µg/L | 5 | 1.92 | 1 | 0.385 |
| Styrene | U | <2.06 | <5.00 | <2.06 | µg/L | 5 | 2.06 | 1 | 0.413 |
| o-Xylene | U | <0.575 | <5.00 | <0.575 | µg/L | 5 | 0.575 | 1 | 0.115 |
| 1,1,2,2-Tetrachloroethane | U | <3.83 | <5.00 | <3.83 | µg/L | 5 | 3.83 | 1 | 0.766 |
| 2-Chlorotoluene | U | <0.660 | <5.00 | <0.660 | µg/L | 5 | 0.660 | 1 | 0.132 |
| 1,2,3-Trichloropropane | U | <3.00 | <5.00 | <3.00 | µg/L | 5 | 3.00 | 1 | 0.599 |
| Isopropylbenzene | J | 2.33 | <5.00 | <0.725 | µg/L | 5 | 0.725 | 1 | 0.145 |
| Bromobenzene | U | <1.33 | <5.00 | <1.33 | µg/L | 5 | 1.33 | 1 | 0.266 |
| n-Propylbenzene | U | <0.680 | <5.00 | <0.680 | µg/L | 5 | 0.680 | 1 | 0.136 |
| 1,3,5-Trimethylbenzene | U | <0.620 | <5.00 | <0.620 | µg/L | 5 | 0.620 | 1 | 0.124 |
| tert-Butylbenzene | U | <0.680 | <5.00 | <0.680 | µg/L | 5 | 0.680 | 1 | 0.136 |
| 1,2,4-Trimethylbenzene | U | <0.570 | <5.00 | <0.570 | µg/L | 5 | 0.570 | 1 | 0.114 |
| 1,4-Dichlorobenzene (para) | U | <1.15 | <5.00 | <1.15 | µg/L | 5 | 1.15 | 1 | 0.23 |
| sec-Butylbenzene | J | 3.96 | <5.00 | <2.26 | µg/L | 5 | 2.26 | 1 | 0.452 |
| 1,3-Dichlorobenzene (meta) | U | <1.07 | <5.00 | <1.07 | µg/L | 5 | 1.07 | 1 | 0.214 |
| p-Isopropyltoluene | U | <0.630 | <5.00 | <0.630 | µg/L | 5 | 0.630 | 1 | 0.126 |
| 4-Chlorotoluene | U | <0.805 | <5.00 | <0.805 | µg/L | 5 | 0.805 | 1 | 0.161 |
| 1,2-Dichlorobenzene (ortho) | U | <1.78 | <5.00 | <1.78 | µg/L | 5 | 1.78 | 1 | 0.357 |
| n-Butylbenzene | U | <0.625 | <5.00 | 0.815 | µg/L | 5 | 0.625 | 1 | 0.125 |
| 1,2-Dibromo-3-chloropropane | U | <4.88 | <25.0 | <4.88 | µg/L | 5 | 4.88 | 5 | 0.977 |
| 1,2,3-Trichlorobenzene | U | <2.00 | <25.0 | <2.00 | µg/L | 5 | 2.00 | 5 | 0.4 |
| 1,2,4-Trichlorobenzene | U | <1.14 | <25.0 | <1.14 | µg/L | 5 | 1.14 | 5 | 0.227 |
| Naphthalene | U | <3.36 | <25.0 | <3.36 | µg/L | 5 | 3.36 | 5 | 0.672 |
| Hexachlorobutadiene | U | <0.990 | <25.0 | 2.36 | µg/L | 5 | 0.990 | 5 | 0.198 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 251 | µg/L | 5 | 250 | 100 | 70 - 130 |
| Toluene-d8 | | 257 | µg/L | 5 | 250 | 103 | 70 - 130 |
| 4-Bromofluorobenzene (4-BFB) | | 256 | µg/L | 5 | 250 | 102 | 70 - 130 |

Sample: 209317 - HLSF-0154-TB-0909-002

Laboratory: Lubbock

Analysis: Volatiles

QC Batch: 63556

Prep Batch: 54254

Analytical Method: S 8260B

Date Analyzed: 2009-09-14

Sample Preparation: 2009-09-14

Prep Method: S 5030B

Analyzed By: JG

Prepared By: JG

continued ...

²⁴Concentration biased low.

sample 209317 continued ...

| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
|------------------------------------|------------------|------------------------|------------------------|---------------------------|-------|----------|--------|---------------------|---------------------|
| Parameter | Flag | SDL Based Result | MQL Based Result | Method Blank Result | Units | Dilution | SDL | MQL (Unadjusted) | MDL (Unadjusted) |
| Bromochloromethane | U | <0.169 | <1.00 | <0.169 | µg/L | 1 | 0.169 | 1 | 0.169 |
| Dichlorodifluoromethane | U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Chloromethane (methyl chloride) | U | <0.164 | <1.00 | <0.164 | µg/L | 1 | 0.164 | 1 | 0.164 |
| Vinyl Chloride | U | <0.110 | <1.00 | <0.110 | µg/L | 1 | 0.110 | 1 | 0.11 |
| Bromomethane (methyl bromide) | U | <0.540 | <5.00 | <0.540 | µg/L | 1 | 0.540 | 5 | 0.54 |
| Chloroethane | U | <0.232 | <1.00 | <0.232 | µg/L | 1 | 0.232 | 1 | 0.232 |
| Trichlorofluoromethane | U | <0.149 | <1.00 | <0.149 | µg/L | 1 | 0.149 | 1 | 0.149 |
| Acetone | 25 JB | 2.17 | <10.0 | 2.05 | µg/L | 1 | 0.893 | 10 | 0.893 |
| Iodomethane (methyl iodide) | U | <0.104 | <5.00 | <0.104 | µg/L | 1 | 0.104 | 5 | 0.104 |
| Carbon Disulfide | U | <0.120 | <1.00 | <0.120 | µg/L | 1 | 0.120 | 1 | 0.12 |
| Acrylonitrile | U | <0.189 | <1.00 | <0.189 | µg/L | 1 | 0.189 | 1 | 0.189 |
| 2-Butanone (MEK) | 26 JB | 0.566 | <5.00 | 0.596 | µg/L | 1 | 0.218 | 5 | 0.218 |
| 4-Methyl-2-pentanone (MIBK) | U | <0.554 | <5.00 | <0.554 | µg/L | 1 | 0.554 | 5 | 0.554 |
| 2-Hexanone | U | <0.863 | <5.00 | <0.863 | µg/L | 1 | 0.863 | 5 | 0.863 |
| trans 1,4-Dichloro-2-butene | U | <0.607 | <10.0 | <0.607 | µg/L | 1 | 0.607 | 10 | 0.607 |
| 1,1-Dichloroethene | U | <0.102 | <1.00 | <0.102 | µg/L | 1 | 0.102 | 1 | 0.102 |
| Methylene chloride | U | <0.270 | <5.00 | <0.270 | µg/L | 1 | 0.270 | 5 | 0.27 |
| MTBE | U | <0.122 | <1.00 | <0.122 | µg/L | 1 | 0.122 | 1 | 0.122 |
| trans-1,2-Dichloroethene | U | <0.168 | <1.00 | <0.168 | µg/L | 1 | 0.168 | 1 | 0.168 |
| 1,1-Dichloroethane | U | <0.107 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| cis-1,2-Dichloroethene | U | <0.151 | <1.00 | <0.151 | µg/L | 1 | 0.151 | 1 | 0.151 |
| 2,2-Dichloropropane | U | <0.167 | <1.00 | <0.167 | µg/L | 1 | 0.167 | 1 | 0.167 |
| 1,2-Dichloroethane (EDC) | U | <0.0890 | <1.00 | <0.0890 | µg/L | 1 | 0.0890 | 1 | 0.089 |
| Chloroform | U | <0.121 | <1.00 | <0.121 | µg/L | 1 | 0.121 | 1 | 0.121 |
| 1,1,1-Trichloroethane | U | <0.170 | <1.00 | <0.170 | µg/L | 1 | 0.170 | 1 | 0.17 |
| 1,1-Dichloropropene | U | <0.156 | <1.00 | <0.156 | µg/L | 1 | 0.156 | 1 | 0.156 |
| Benzene | U | <0.113 | <1.00 | <0.113 | µg/L | 1 | 0.113 | 1 | 0.113 |
| Carbon Tetrachloride | U | <0.0930 | <1.00 | <0.0930 | µg/L | 1 | 0.0930 | 1 | 0.093 |
| 1,2-Dichloropropane | U | <0.129 | <1.00 | <0.129 | µg/L | 1 | 0.129 | 1 | 0.129 |
| Trichloroethene (TCE) | U J | <0.160 | <1.00 | <0.160 | µg/L | 1 | 0.160 | 1 | 0.16 |
| Dibromomethane (methylene bromide) | U | <0.137 | <1.00 | <0.137 | µg/L | 1 | 0.137 | 1 | 0.137 |
| Bromodichloromethane | U | <0.107 | <1.00 | <0.107 | µg/L | 1 | 0.107 | 1 | 0.107 |
| 2-Chloroethyl vinyl ether | U | <0.257 | <5.00 | <0.257 | µg/L | 1 | 0.257 | 5 | 0.257 |
| cis-1,3-Dichloropropene | U | <0.109 | <1.00 | <0.109 | µg/L | 1 | 0.109 | 1 | 0.109 |
| trans-1,3-Dichloropropene | U | <0.181 | <1.00 | <0.181 | µg/L | 1 | 0.181 | 1 | 0.181 |
| Toluene | U | <0.117 | <1.00 | <0.117 | µg/L | 1 | 0.117 | 1 | 0.117 |
| 1,1,2-Trichloroethane | U | <0.143 | <1.00 | <0.143 | µg/L | 1 | 0.143 | 1 | 0.143 |
| 1,3-Dichloropropane | U | <0.191 | <1.00 | <0.191 | µg/L | 1 | 0.191 | 1 | 0.191 |
| Dibromochloromethane | U | <0.118 | <1.00 | <0.118 | µg/L | 1 | 0.118 | 1 | 0.118 |

continued ...

²⁵ Concentration biased low.²⁶ Concentration biased low.

sample 209317 continued ...

| Parameter | Flag | SDL | MQL | Method | Units | Dilution | SDL | MQL | MDL |
|-----------------------------|-----------------|--------------|--------------|--------------|-------|----------|-------|--------------|--------------|
| | | Based Result | Based Result | Blank Result | | | | (Unadjusted) | (Unadjusted) |
| 1,2-Dibromoethane (EDB) | ^U | <0.131 | <1.00 | <0.131 | µg/L | 1 | 0.131 | 1 | 0.131 |
| Tetrachloroethene (PCE) | ^{27 U} | <0.353 | <1.00 | <0.353 | µg/L | 1 | 0.353 | 1 | 0.353 |
| Chlorobenzene | ^U | <0.135 | <1.00 | <0.135 | µg/L | 1 | 0.135 | 1 | 0.135 |
| 1,1,1,2-Tetrachloroethane | ^U | <0.263 | <1.00 | <0.263 | µg/L | 1 | 0.263 | 1 | 0.263 |
| Ethylbenzene | ^U | <0.174 | <1.00 | <0.174 | µg/L | 1 | 0.174 | 1 | 0.174 |
| m,p-Xylene | ^U | <0.220 | <1.00 | <0.220 | µg/L | 1 | 0.220 | 1 | 0.22 |
| Bromoform | ^U | <0.385 | <1.00 | <0.385 | µg/L | 1 | 0.385 | 1 | 0.385 |
| Styrene | ^U | <0.413 | <1.00 | <0.413 | µg/L | 1 | 0.413 | 1 | 0.413 |
| o-Xylene | ^U | <0.115 | <1.00 | <0.115 | µg/L | 1 | 0.115 | 1 | 0.115 |
| 1,1,2,2-Tetrachloroethane | ^U | <0.766 | <1.00 | <0.766 | µg/L | 1 | 0.766 | 1 | 0.766 |
| 2-Chlorotoluene | ^U | <0.132 | <1.00 | <0.132 | µg/L | 1 | 0.132 | 1 | 0.132 |
| 1,2,3-Trichloropropane | ^U | <0.599 | <1.00 | <0.599 | µg/L | 1 | 0.599 | 1 | 0.599 |
| Isopropylbenzene | ^U | <0.145 | <1.00 | <0.145 | µg/L | 1 | 0.145 | 1 | 0.145 |
| Bromobenzene | ^U | <0.266 | <1.00 | <0.266 | µg/L | 1 | 0.266 | 1 | 0.266 |
| n-Propylbenzene | ^U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,3,5-Trimethylbenzene | ^U | <0.124 | <1.00 | <0.124 | µg/L | 1 | 0.124 | 1 | 0.124 |
| tert-Butylbenzene | ^U | <0.136 | <1.00 | <0.136 | µg/L | 1 | 0.136 | 1 | 0.136 |
| 1,2,4-Trimethylbenzene | ^{28 J} | 0.269 | <1.00 | <0.114 | µg/L | 1 | 0.114 | 1 | 0.114 |
| 1,4-Dichlorobenzene (para) | ^U | <0.230 | <1.00 | <0.230 | µg/L | 1 | 0.230 | 1 | 0.23 |
| sec-Butylbenzene | ^U | <0.452 | <1.00 | <0.452 | µg/L | 1 | 0.452 | 1 | 0.452 |
| 1,3-Dichlorobenzene (meta) | ^U | <0.214 | <1.00 | <0.214 | µg/L | 1 | 0.214 | 1 | 0.214 |
| p-Isopropyltoluene | ^U | <0.126 | <1.00 | <0.126 | µg/L | 1 | 0.126 | 1 | 0.126 |
| 4-Chlorotoluene | ^U | <0.161 | <1.00 | <0.161 | µg/L | 1 | 0.161 | 1 | 0.161 |
| 1,2-Dichlorobenzene (ortho) | ^U | <0.357 | <1.00 | <0.357 | µg/L | 1 | 0.357 | 1 | 0.357 |
| n-Butylbenzene | ^U | <0.125 | <1.00 | 0.163 | µg/L | 1 | 0.125 | 1 | 0.125 |
| 1,2-Dibromo-3-chloropropane | ^U | <0.977 | <5.00 | <0.977 | µg/L | 1 | 0.977 | 5 | 0.977 |
| 1,2,3-Trichlorobenzene | ^U | <0.400 | <5.00 | <0.400 | µg/L | 1 | 0.400 | 5 | 0.4 |
| 1,2,4-Trichlorobenzene | ^U | <0.227 | <5.00 | <0.227 | µg/L | 1 | 0.227 | 5 | 0.227 |
| Naphthalene | ^U | <0.672 | <5.00 | <0.672 | µg/L | 1 | 0.672 | 5 | 0.672 |
| Hexachlorobutadiene | ^U | <0.198 | <5.00 | 0.473 | µg/L | 1 | 0.198 | 5 | 0.198 |

| Surrogate | Flag | Result | Units | Dilution | Spike Amount | Percent Recovery | Recovery Limits |
|------------------------------|------|--------|-------|----------|--------------|------------------|-----------------|
| Dibromofluoromethane | | 48.6 | µg/L | 1 | 50.0 | 97 | 70 - 130 |
| Toluene-d8 | | 50.1 | µg/L | 1 | 50.0 | 100 | 70 - 130 |
| 4-Bromofluorobenzene (4-BFB) | | 49.8 | µg/L | 1 | 50.0 | 100 | 70 - 130 |

Method Blank (1)QC Batch: 63556
Prep Batch: 54254Date Analyzed: 2009-09-14
QC Preparation: 2009-09-14Analyzed By: JG
Prepared By: JG²⁷Concentration biased low.²⁸Concentration biased low.

705 S. TELEGRAPH ST. F-201
 LEB CITIES, MO 63011
 616-602-1526
 616-602-1501

CHAIN OF CUSTODY RECORD

PROJECT NO.

PROJECT NAME

HEISTE Semi-Annual Ground Water

SAMPLES TO BE ANALYZED

Bridget Davis

ANALYSIS REQUESTED

REMARKS

PAGE 1 OF 1

0002

| DATE | TIME | SAMPLE ID | MATERIAL | LAB NO. | NO. OF CONTAINERS | ANALYSIS REQUESTED | REMARKS |
|--------|------|------------------------|--------------|---------|-------------------|--------------------|-------------------------|
| 9-1-09 | 1335 | HLSE-D154-DRU-016-0909 | Water | 209311 | 3 | X | |
| 9-2-09 | 1115 | HLSE-0154-HCF-081-0909 | Water | 312 | 3 | X | |
| 9-1-09 | 1033 | HLSE-0154-DRU-005-0909 | Water | 313 | 3 | X | |
| 9-1-09 | 1030 | HLSE-0143-HMW-013-0909 | Water | 314 | 3 | X | |
| 9-1-09 | 1010 | HLSE-0143-HMW-036-0909 | Water | 315 | 3 | X | |
| 9-1-09 | 1335 | HLSE-0154-DRU-016-0909 | ms/msd Water | 209311 | 3 | X | Extra sample for ms/msd |
| 9-2-09 | 1338 | HLSE-0154-DRU-004-0909 | Water | 316 | 3 | X | |
| 9-2-09 | 1338 | HLSE-0154-DRU-002-0909 | Water | 317 | 3 | X | trip Blank |

| PROJECT INFORMATION | | SAMPLES RECEIVED | | 1. RECEIVED BY (SIGNATURE) | | 2. RECEIVED BY (SIGNATURE) | | 3. RECEIVED BY (SIGNATURE) | |
|---------------------|---------------|---------------------------|--|----------------------------|--|----------------------------|--|----------------------------|--|
| PROJECT MANAGER | | TOTAL NO. OF CONTAINERS | | (PRINTED NAME) | | (PRINTED NAME) | | (PRINTED NAME) | |
| FIELD DATE | | CHAIN OF CUSTODY SEALS | | RECEIVED BY (SIGNATURE) | | RECEIVED BY (SIGNATURE) | | RECEIVED BY (SIGNATURE) | |
| SHIPPING ID NO. | | GOOD CONDITION MAINTAINED | | (PRINTED NAME) | | (PRINTED NAME) | | (PRINTED NAME) | |
| TRAILER NO. | 7099080805084 | CONFORMS TO RECORD | | (PRINTED NAME) | | (PRINTED NAME) | | (PRINTED NAME) | |
| VEHICLE | Fed EX | | | | | | | | |

PLEASE SEE ATTACHED ANALYSIS LIST FOR DETAILS